

April 16, 2014

Ms. Kimberly Tisa
United States Environmental Protection Agency
5 Post Office Square, Suite 100
Mail Code: OSRR07-2
Boston, MA 02109-3912

RE: Application for Risk-Based Soil Cleanup Plan
Maine Energy, Biddeford, Maine

Dear Ms. Tisa:

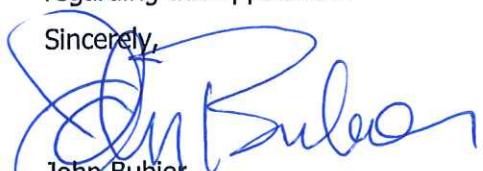
Maine Energy is requesting approval of a risk-based Polychlorinated Biphenyl cleanup application from U.S. EPA Region 1 (EPA) under 40 CFR 761.61(c). The attached risk-based application presents sampling results, a risk analysis, and a proposed remediation plan to remove and dispose of approximately 1,082 cubic yards of demolition debris and fill containing PCBs over 50 parts per million (ppm) and approximately 1,000 cubic yards of demolition debris and fill containing PCBs greater than 25 ppm but less than 50 ppm. The application also presents details on a pending environmental covenant to permanently restrict the area of concern to certain land uses.

The enclosed risk-based application is supported by analytical data from over 150 soil samples collected within and adjacent to the area of concern. The risk analysis presented in Section 5 provides support for the conclusion that collectively, the soil remedy and the cap proposed to be installed, are protective of human health and the environment.

Our tentative schedule to implement the proposed Plan is during early summer of this year. Maine Energy will initiate the activities described in the application upon approval by EPA.

Please contact me if we can provide additional information or if you have any questions regarding this application.

Sincerely,



John Bubier
City of Biddeford



Brian Oliver
Maine Energy Recovery Company, LP

CC: John K. Cressey, Summit
Brian Phinney, City of Biddeford
Ken Robbins, Maine Energy Recovery Company, LP
Nicholas Hodgkins, Maine DEP



**APPLICATION FOR
40 CFR 761.61(c) RISK-BASED CLEANUP OF SOIL
MAINE ENERGY
BIDDEFORD, MAINE**

Prepared For:
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April 16, 2014
Project # 12-3259.1

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1.0 INTRODUCTION

This risk-based application is for cleanup of polychlorinated biphenyl (PCB) impacted demolition debris and fill at a former boilerhouse on the eastern portion of the former Maine Energy property in Biddeford, Maine. The City of Biddeford completed its purchase of the property in November 2012. A Site Locus Map is included as Figure 1. This application includes:

1. Results for characterization and delineation of PCB impacted demolition debris and fill (completed).
2. A Plan for removal and off-site disposal of demolition debris and fill containing PCBs greater than 25 parts per million (ppm) and backfilling with clean fill (planned).
3. A description of a multi-layer capping system being proposed for installation to prevent exposure to remaining (underlying) PCB impacted fill.
4. A proposal to execute a deed covenant restricting certain land uses on the area of concern.
5. A verification sampling plan following soil removal (planned).

This risk-based cleanup application is consistent with the risk-based requirements of 761.61(c) with the exception that a modified characterization sampling approach was completed under Subpart N of the CFR (cleanup site characterization sampling). Justification for the modified characterization sampling is presented in Sections 3 and 4, respectively.

1.1 Background

The Maine Energy Recovery Company waste to energy (incinerator) facility was constructed on an 8.4 acre parcel of property in downtown Biddeford, Maine in 1984. The parcel is bounded by the Saco River on the east, Pan Am Railway to the north, Lincoln Street to the west and Pearl Street to the south. The facility consisted of several industrial and office buildings. In 2012, Maine Energy and the City of Biddeford reached an agreement to close the facility and enter into an agreement to allow the City to purchase the property.

According to a Phase I Environmental Site Assessment (ESA) completed on the property in 2012

The earliest information for the Site was an 1885 Sanborn Fire Insurance map that indicated the Site and surrounding area had been heavily developed as a downtown area with mill buildings, other industrial buildings, and numerous small commercial buildings and dwellings. Primary uses of the area included foundry operations and machining, textile manufacturing, warehousing, and trucking or automotive services.

In the fall of 2012, a Phase I and Phase II Environmental Site Assessment was performed on behalf of the City of Biddeford at the Maine Energy facility. Site investigations documented PCB Aroclor 1260 as being present within demolition debris and fill along the eastern boundary of the property (area of concern) consistent with the location of a former boilerhouse. The Sanborn Fire Insurance Maps (Sanborn) from 1932 and 1947 and aerial photographs from between 1956 to 1975 depict the boilerhouse as being located on this parcel; however, the 1976 Sanborn Map and aerial photographs from 1986, 1990, 1998, and depict the parcel as vacant. See Appendix A for copies of these maps and aerial photographs. Maine Energy structures were not built on the area of concern and the property remains vacant with the exception of three electrical transmission poles that were installed to transfer electricity from the power generating facility located to the west of the area of concern to the electrical grid on

the opposite (east) side of the Saco River. These poles were subsequently cut off at ground surface as a portion of the demolition of the Maine Energy facility in late 2013.

Property to the north owned by the Boston & Maine Railroad (Pan Am) is an active elevated rail line (Amtrak Downeaster commuter) which has 5 round trip commuter trains per day seven days a week serving Boston. A retaining wall for the rail line is located on/near the boundary of the Maine Energy parcel and the Pan Am Right of Way (ROW) and is the northern boundary of the area of concern. Due to stability concerns for the retaining wall and rail line, a Geotechnical Evaluation of the retaining wall and rail line stability was performed. Results of the Geotechnical Evaluation indicate that excavations near the retaining wall should be minimized to avoid stability concerns. As such, the cleanup plan has incorporated measures to minimize excavation in the vicinity of the retaining wall, while achieving a cleanup that is protective of human health and the environment.

An evaluation of engineering options to support the retaining wall show that costs are extremely high, stability concerns cannot be completely eliminated and the reduction in risk by excavating additional PCB impacted fill at depth is minimal. Therefore, the cleanup plan proposes that PCB impacted fill at concentrations below 25 parts per million (ppm) be left on site, approximately 4 feet below grade, and covered with an enhanced capping system to eliminate exposure pathways and minimize risk associated with PCB impacted Fill remaining at the Site.

1.2 Application

This application is organized in the following sections:

Section 2 – Remedial Approach and Objectives: describes the general approach to remediation, occupancy levels, restrictions, cleanup goals, and corresponding risk levels.

Section 3 – Sampling Approach and Characterization Data: describes the various phases of soil sampling and analysis investigations conducted between 2012 and 2013.

Section 4 – Soil Remediation Plan describes the removal, off-site disposal, and verification sampling of soil from the areas containing PCBs greater than 25 ppm.

Section 5 –Risk Analysis: describes the risk evaluation to potential human receptors from contact with PCBs in fill beneath a proposed capping system.

Section 6 –References: lists the references used to support the application

Tables, figures, and appendices follow the text.

2.0 REMEDIAL APPROACH AND OBJECTIVES

The overall goal of the Maine Energy remediation project is to protect human health and the environment by remediating and managing PCB Aroclor 1260 impacted demolition debris and fill at the Maine Energy parcel. To achieve this goal, a proposed cleanup plan has been developed consistent with the Toxic Substances Control ACT (TSCA) and Maine Department of Environmental Protection (MEDEP) Voluntary Response Action Program (VRAP) standards.

2.1 Facility Boundaries and Ownership

The portion of the former Maine Energy property that this plan is designed to remediate is an approximately 0.45 acre parcel of land bound by the Pan Am Railroad property to the north, the Saco River to the south and east, and the remainder of the Maine Energy facility to the west. Maine Energy has owned the facility since construction commenced. After approximately 2 years of construction, the facility began operation in 1987. The City of Biddeford has purchased the parcel from Maine Energy, although Maine Energy will be responsible for implementing this cleanup plan.

2.2 Occupancy and Restrictions

The current reuse of this portion of the property is planned to be a public park with greater than 6.7 hours per week of access for Park Users, the defined period for "high occupancy" use under the PCB definitions (40 CFR 761.3). The remainder of the property is being considered for potential residential and commercial uses.

The City of Biddeford has agreed to execute an Environmental Deed Covenant under the Uniform Environmental Covenants Act (UECA) Maine Revised Statutes 38 MRSA 3001 to restrict land use to a park or similar public open space with conditions prohibiting the disturbance of subsurface soil and fill below 4 feet from ground surface. Both the property owner, City of Biddeford, current tenant Maine Energy, as well as the approving agency, Maine DEP, will be holders of the environmental covenant. The environmental covenant will be recorded at the York County Registry of Deeds.

2.3 PCB Waste Classification and Cleanup Levels

Demolition debris and fill that contains PCB 1260 is considered PCB remediation waste under 761.3. Accordingly, management and disposal of PCB remediation waste is based on the "as found" PCB concentration of individual samples collected in situ. Impacts appear to be associated with the former boilerhouse and therefore, the remediation of the Site is correlated with the location of this building. As referenced on Figure 4, impacts observed in analytical results are highest in the center and eastern portions of this former structure. Concentrations of PCBs decline rapidly outside of the building footprint and to the west. Areas 1, 2, and 3 as shown on Figure 5 contain the highest concentrations of PCBs and will be managed as hazardous (≥ 50 ppm) waste while Areas 4 and 5 contain lower concentrations of PCBs and will be managed as non-hazardous (< 50 ppm) waste.

The proposed cleanup standard is a goal of a minimum removal of 4 feet of soil but additional excavation will be completed to obtain the cleanup goal of less than 25 ppm.

Proposed cleanup levels for PCB remediation waste according to location of the samples are listed below.

1. Areas 1 and 2—Based upon analytical results (summarized in Section 3.1), demolition debris and fill containing greater than 25 ppm in individual soil core samples will be

excavated as structurally feasible (summarized in Section 4.2.1.1) and replaced with clean fill and a cap system. The excavated soil will be managed as PCB remediation waste (note that analytical results for some samples from these areas reported PCB concentrations greater than 50 ppm) and be subject to disposal requirements under TSCA and disposed as hazardous waste at a facility licensed to accept this waste.

2. Area 3 – Based upon analytical results, demolition debris and fill containing greater than 50 ppm in individual soil core samples will be excavated to a depth of 4 feet below ground surface and removed as TSCA Remediation Waste and disposed as hazardous waste at a facility licensed to accept this waste. Once the fill has been removed, a protective cap as described in Section 4 will be installed in this area.
3. Areas 4 – Based upon analytical results, PCB impacted fill with concentration between 1 and 25 ppm are present. Fill will be excavated to a depth of 4 feet below grade and replaced with a protective cap as described in Section 4. The excavated soil will be managed as <50 ppm PCB waste and disposed at Juniper Ridge Landfill in Old Town, Maine.
4. Area 5 – Based upon analytical results, PCB impacted fill with concentration between 1 and 2.24 ppm are present. Fill will be excavated to a depth of 2 feet below grade (but will extend to 4 feet if the retaining wall footer elevation allows for it) and replaced with a protective cap as described in Section 4. The excavated soil will be managed as <50 ppm PCB waste and disposed at Juniper Ridge Landfill in Old Town, Maine.
5. Area 6 – Based upon analytical results, PCB impacted demolition debris and fill containing greater than 25 ppm in individual soil core samples will be excavated as structurally feasible and replaced with clean fill and a cap system. The excavated soil will be managed as PCB remediation waste and be subject to disposal requirements under TSCA and disposed as hazardous waste at a facility licensed to accept this waste.

This area will be sloped away from the retaining wall with excavation beginning at ground surface and continuing to the top of the retaining wall footer then sloping at a ratio of 1 foot vertical to 1.5 feet horizontal (i.e. if the bottom of the footer were at 4 feet below ground surface, excavation would commence to 4 feet then slope to 5 feet at a distance of 1.5 feet and continue in the manner until 10 feet (or the 25 ppm remediation goal is reached).

3.0 SAMPLING APPROACH AND CHARACTERIZATION DATA

Sampling of fill and soil as bulk PCB remediation waste was designed to adequately characterize the vertical and horizontal extent of impact, acknowledging the physical features of the site (retaining walls, transmission poles), current and planned occupancy, and the presumed type of release. There is no known date or point of release; rather, the past presence of the boilerhouse and potential PCB-containing materials within the building are believed to have been released and distributed proximal to the former boilerhouse. The PCB that is found on the property, Aroclor 1260, is solid at room temperature and almost insoluble in water. As a result, the mobility of Aroclor 1260 in the environment is limited and typically would be associated with the physical transport or disturbance of soil particles to which it is bound. The target area for delineation and characterization of PCB impact was the area proximal to the former boilerhouse. Topography of the area is relatively flat with approximately four feet of topographic relief across the former boilerhouse area, sloping from west to east. Approximately half of the area of concern is asphalt paved (former trailer parking area) and the other half is grassy/meadow area. The area of concern is bounded on the north and east sides by retaining walls (concrete wall to the north toward the rail line and granite blocks to the east). The retaining walls extend 1-3 feet above grade and act as a physical barrier for movement or transport of soil/sediment beyond the retaining wall.

Materials observed in borings included loam, granular fill (sand and gravel), brick, concrete, coal ash/clinker, cobbles and boulders. Based on analytical results from multiple investigations conducted in 2012 and 2013, PCB concentrations were found to be variable throughout the demolition debris and fill matrix and not correlated to a specific depth or type of fill/debris. These data suggest that the PCB impacts were not associated with a discrete spill or release, but were likely distributed during demolition of the former boilerhouse.

Over 150 soil samples were collected from within and adjacent to the area of concern for laboratory analysis since October 2012. Soil sampling methods consisted of direct-push technology and standard penetration test technology. Soil core diameters ranged from 1.5 to 4 inches. Select samples were submitted to Absolute Resource Associates in Portsmouth, New Hampshire for sampling completed in October 2012 and January 2013 and to Analytics Environmental Laboratory in Portsmouth, New Hampshire for sampling completed in February 2013. All samples were submitted for laboratory analysis of total PCBs using EPA Method 8082 (soxhlet extraction). Sampling locations are shown on Figure 3.

It should be noted that the original Phase II ESA investigation of the parcel included the analysis of 69 samples using immunoassay test kits following USEPA Method 4020. These initial sample results were used as a screening tool to preliminarily assess the magnitude and extent of impacts.

3.1 Soil Sampling on Site

3.1.1 Phase II ESA – October 2012

On October 2nd and 3rd, 2012, Credere Associates, LLC (Credere) completed a Phase II Environmental Site Assessment (ESA) on the Maine Energy property on behalf of the City of Biddeford (City) in conformance with the ASTM International (ASTM) E1903-11 Standard Practice for Environmental Site Assessments: Phase II Environmental Site Assessment Process.

This Phase II ESA assessed environmental media at the Site to confirm or dismiss thirteen (13) recognized environmental conditions (RECs) identified during a Phase I ESA (dated September 18, 2012) that was completed by Credere.

A summary of the work tasks completed during the Phase II ESA are included below:

1. A Site specific Health and Safety Plan (HASP) was developed for investigatory personnel.
2. The investigation locations were approved by Maine Energy and pre-marked by Credere. The Site was cleared for underground utilities by DigSafe, and each individual soil boring location was pre-cleared by DigSmart of Maine using electronic locating equipment and ground penetrating radar (GPR).
3. Fifteen (15) soil borings were advanced throughout the Site to facilitate the collection of soil samples for visual logging, field screening, and laboratory analysis.
4. Six (6) of the soil borings were completed as groundwater monitoring wells and ground water samples were collected for laboratory analysis. Note that up to nine (9) wells were proposed to be installed at the Site, but a saturated zone in the overburden soil was not observed in some areas due to shallow bedrock. Each monitoring well was subsequently surveyed to allow for the calculation of approximate groundwater flow directions at the Site.
5. Five (5) surficial soil samples were collected from the eastern portion of the Site where ash was observed at the ground surface.
6. Five (5) concrete floor samples were collected from within the Site building(s) for laboratory analysis.
7. One (1) sample was collected from the interior brick lining of Maine Energy's exhaust stack for laboratory analysis.

The Phase II investigation included the collection of seventeen soil samples for analysis of Polychlorinated Biphenyls (PCBs). Of the seventeen samples collected, 15 reported PCBs as Not Detected (ND), two samples reported PCB concentrations above the laboratory detection limit but only one sample, CA-SB-5, reported PCB concentrations above the US EPA TSCA unit's notification limit of 1 part per million (ppm). Sample CA-SB-5 was collected from a boring located in the area of the former boilerhouse.

Analytical results of Sample CA-SS-4 reported PCB concentrations above the laboratory detection limit (0.04 ppm), but below the TSCA notification limit of 1 ppm. CA-SS-4 was collected near a 20,000 gallon underground storage tank (UST) along the northern boundary of the property. Additional surface soil samples (CA-SS-1, CA-SS-2, CA-SS-3, and CA-SS-5) were collected at locations across the eastern portion of the property and were reported as non-detect for PCBs.

Based upon Phase II analytical results, Credere recommended a supplemental Phase II ESA to further investigate the potential for PCB impacts on the eastern portion of the property. See Figure 2 for a copy of the Phase II ESA site plan showing sampling locations and Table 1 for analytical results for PCB samples collected during the Phase II ESA.

3.1.2 Supplemental Phase II ESA – October 2012

Following receipt and interpretation of analytical data from the Phase II ESA, a supplemental Phase II ESA was completed by Credere on October 25, 2012. The Supplemental Phase II ESA included the advancement of 19 borings in the area of the former boilerhouse. Soil samples were collected from each boring and screened for PCBs by Cape Technologies of South Portland, Maine using Immunoassay test kits. Four samples analyzed by Cape Technologies

were split and submitted to Absolute Resource Associates (Absolute) of Portsmouth, New Hampshire for analysis of PCBs using soxhlet extraction by US EPA Method 8082. Results of two of the four samples analyzed by Absolute reported PCB concentrations above the 1 ppm TSCA notification limit and were generally consistent with the immunoassay test kit results. The Immunoassay testing methodology utilized by Cape Technologies is considered a screening methodology by EPA, and is not appropriate for use in developing a soil removal Work Plan without extensive verification and validation testing to demonstrate the precision of the methodology. Nonetheless, the investigation completed by Credere suggested that PCBs were present in fill near the former Boiler House. See Figure 3 for locations of the investigation, Table 2 for immunoassay test results, and Table 3 for analytical results collected for correlation with immunoassay results.

3.1.3 Characterization Investigation – January 2013

Based on the Phase II ESA investigations, Summit completed a characterization and delineation investigation in January 2013 to collect additional samples for analysis of PCB in the area of concern. Summit advanced twelve soil borings (SE-SB-201 through SE-SB-212) between monitoring well CA-MW-10 and the section of pavement where trailers were formerly parked. See Figure 4 for a boring location map.

Summit collected soil samples from visually distinctive intervals and submitted them to Absolute for analysis of PCBs by EPA Method 8082 using soxhlet extraction. Analytical results reported PCB concentrations ranging from Non-Detect ND (i.e., less 0.03 ppm) to 3,500 ppm. PCB impacts were generally associated with demolition debris, but not necessarily with specific depth or type of debris (i.e., brick concrete, granular soil, ash). This round of sampling narrowed the extent of impacts. See Table 4 for analytical results.

3.1.4 Supplemental Characterization Investigation – February 2013

Summit discussed the October 2012 and January 2013 results with you on January 15, 2013 to ascertain the next steps that should be taken to collect additional data prior to completion of this application. Based on this discussion, a horizontal grid of 20 feet by 20 feet and soil samples collected every two feet vertically was considered to be appropriate. On February 19th and 20th, 2013 Summit advanced twenty-four (24) additional soil borings (SE-SB-213 through SE-SB-236) and sampled these soil borings from each vertical 2 foot interval for PCBs. Soil borings were advanced to refusal or native soil. Materials observed in these borings were loam, sand and gravel fill, brick, concrete, ash, and silty-clay (native). Soil samples were submitted to Analytics for analysis of PCBs by EPA Method 8082 using soxhlet extraction. Analytical results reported PCB concentrations ranging from ND to 9,660 ppm. PCB impacts could not be associated with a specific material or soil interval, however, impacts are generally coincident with the footprint of the former Boiler House. See Table 5 for analytical results.

3.2 Characterization Summary

Demolition debris and fill in the vicinity of the former boilerhouse was sampled at over 50 locations using a combination of screening methodology, focused sampling (e.g. specific materials within and outside of the former boilerhouse footprint) and systematic sampling (e.g. horizontal and vertical grid coordinates). Sampling was conducted over several iterative investigations between October 2012 and February 2013. The results from all sampling assessments reveal the following patterns:

- The highest concentrations are found in the soils located within the footprint of the former boilerhouse;
- PCB concentrations rapidly decrease with depth and distance from the former boilerhouse.
- Concentrations of PCBs in the area of concern range from non-detect to 9,600 ppm; 21 samples out of 152 total samples exceed 50 ppm and 26 samples out of the 152 samples exceed 25 ppm.

The distribution pattern and concentrations of PCBs are consistent with the presence of PCB impacts associated with the boilerhouse and subsequent redistribution resulting from disturbance and/or demolition of the boilerhouse.

4.0 SOIL REMEDIATION PLAN

The proposed soil Remediation Plan consists of removing approximately 1,1,082 cubic yards (1,1,623 Tons) of demolition debris and fill with PCB concentrations greater than 50 ppm. Management and disposal of this excavated material will be as described in section 4.2.3 below. The excavated volume includes all fill characterized as potentially having PCB concentrations greater than 50 ppm. The area of fill containing concentrations greater than 50 ppm is delineated on Figure 5.

Following excavation of PCB impacted fill with concentrations greater than 50 ppm, fill from 0-4 feet below grade will be excavated and disposed off-site as described in Section 4.2.4. It is estimated that approximately 1,000 cubic yards (1,500 Tons) will be removed from the area delineated on Figure 5. PCB concentrations range from ND to less than 15.8 ppm within this excavation area. Due to the variability of PCB concentrations additional segregation of the fill (i.e., some areas may be less than 1 ppm) cannot feasibly be completed. Therefore, a uniform 4 feet will be removed and an enhanced cap system will be installed. The depth of 4 feet below grade was selected based on the following factors:

- A geotechnical evaluation indicated that deeper excavations pose a stability concern for the abutting retaining wall to the north. This wall supports an active rail line (Downeaster Amtrak commuter line).
- A four foot thick enhanced capping system will eliminate a direct contact exposure route for remaining underlying PCB impacted fill. Note that PCB concentrations in fill below the capping system range from ND to less than 25 ppm.
- An Environmental Deed Covenant will be recorded to establish Institutional Control for the cleanup area. The Deed Covenant will restrict land use and activities within the cleanup area such that remaining PCB impacts will not pose a risk to human health or the environment. The covenant will also include long-term maintenance requirements to maintain the integrity of the capping system.
- Four feet is considered well beyond the depth of typical disturbance associated with a park setting (benches, lawn maintenance, shrub and tree planting, walkway construction, etc.), thereby being protective of exposure due to inadvertent activity, even if restricted by Covenant.
- The cost to design and implement an engineering control to stabilize the retaining wall is prohibitive [timber lag system drilled into bedrock and braced]. Remediation cost is estimated to triple in order to excavate an additional 4-8 feet of demolition debris and fill impacted with PCB concentrations less than 25 ppm. See Section 4.2.1.1 below.
- Any Engineered Stabilization Plan would need the approval of Pan Am Railways and Amtrak. This process is unlikely to allow timely implementation of a Remediation Plan.

4.1 Soil Excavation

Based on the results of characterization sampling completed within this area, PCB impacted materials appear to have been reworked and distributed within the foundation of the former boilerhouse. Impacted materials are present within this area at differing depths but appear to be consistent with demolition debris or fill associated with the former boilerhouse.

4.2 Field Procedures

Plans and specs for the soil removal will be prepared and used to select contractors prior to initiating work. All work will be conducted in accordance with a site-specific health and safety

plan (HASP). The following sections describe the general work elements of the soil removal action.

4.2.1 Excavation

Each excavation area will be marked in advance and cleared for utilities. Excavation will be performed by mechanical means. The dimensions and expected removal volumes based upon removal of soil containing levels of PCBs greater than 25 ppm are listed on Figure 5. The total volume removed is anticipated to be approximately 2,082 cubic yards or roughly 3,123 tons.

Due to structural considerations with the railroad retaining wall (See Figure 4), the excavation of all PCB-impacted material is not cost effective. The excavation may additionally be limited by groundwater (observed between 11 and 14 feet below ground surface).

4.2.1.1 Geotechnical Analysis

Summit Geotechnical Services (SGS) completed a geotechnical evaluation of the concrete retaining wall which supports the elevated rail line. The geotechnical recommendations for remediation excavation slopes, dewatering, and potential debris removal are provided as follows:

Concrete Retaining Wall

Portions of excavations are anticipated adjacent to retaining wall structures (gravity walls). Excavations are planned down slope of a concrete retaining wall 13 feet in height located along the northern portion of the site retaining soils located beneath the Pan Am Railroad. The railroad line is located approximately 25 to 45 feet north of the retaining wall.

Exact depth and dimension of the concrete retaining wall footings are not known. Hand probes performed adjacent to the retaining wall suggest a consistent refusal depth of 1 to 1.5 feet below existing grade and may represent the location for the top of retaining wall footing. For excavation stability and to prevent undermining of the retaining wall footing, SGS recommends a minimum excavation slope of 1.5 horizontal to 1 vertical be maintained extending from the toe/base of the footing. Alternatively braced excavations such as underpinning and shoring will be required. Maximum excavation depth within the area of concern should be limited to a maximum depth of 10 feet below existing top of grade and less than 6 feet below the base of footing elevation. Excavation adjacent to the retaining wall should be performed using staged excavations with a maximum length of exposed wall limited to 15 feet or less.

Due to the presence of rubble debris including brick, concrete, and ash it is recommended that remediation excavations adjacent to the retaining walls be inspected by the geotechnical engineer to verify subgrade is in conformance with design parameters and stable to support excavation slopes of 1.5 horizontal to 1 vertical. Using the provided topographic information available, boring exploration data, and measurements obtained during our site visit we estimate the following factors of safety for remediation excavation adjacent to the concrete retaining wall:

Global Slope Stability (Slide 5.0, Bishop Simplified), FS = 1.3

Lateral Base Sliding (Estimated footing width of 7 feet), FS = 1.5

Overturning (Estimated footing width of 7 feet), FS = 1.5

Bearing Capacity (Maximum slope cut of 1.5H to 1V), FS = 3

Excavation Slopes

The subgrade is classified as type C soil in conformance with Occupational Safety and Health Administration (OSHA) excavation guidelines. Excavations should be limited to a maximum side slope of 1.5 horizontal to 1 vertical to a maximum depth of 10 feet. Excavations with a depth greater than its width and a width less than 15 feet are considered trench excavations by OSHA and require conformance to applicable OSHA trench excavations standards. Where excavations are performed below a depth of 4 feet with widths less than 15 feet, OSHA trench standards may apply requiring the use of a trench box or similar excavation support system. Excavations adjacent to the concrete retaining wall should be limited to a length of wall exposure of 15 feet or less.

Where excavation is performed adjacent to existing structures, a series of monitoring points including settlement monitoring points (SMP) and deflection monitoring points (DMP) should be considered to monitor potential movements. Suggested locations for monitoring points include the concrete retaining wall and granite block retaining wall where extended time of open excavation (greater than 24 hours) is anticipated.

The SMP and DMP should be installed and monitored by a Maine Licensed Surveyor or qualified personnel using sufficiently accurate survey equipment. All SMP and DMP should be clearly marked for easy identification. Monitoring points may consist of nails and screws, reinforcing bars, or similar materials with well-defined measurement points. Where available observable cross marks on the top of horizontal structures may also be used. The surface within 3 inches of the point should be cleaned and/or marked to permit easy identification. The monitoring points to measure the elevations of the DMP and SMP should be recorded to 0.01 ft for position of DMP and the horizontal position of SMP measured to 0.1 foot. See Figure 7 for an Excavation Plan from SGS.

Excavation Dewatering

SGS recommends surface drainage be diverted from open excavations to prevent the collection of water during wet periods such as rain or snowmelt. Temporary surface diversions such as ditching, sloping, sand bags or other suitable method to adequately prevent surface flow from entering excavations should be employed prior to excavation.

Excavation below the water table will be avoided when possible to prevent additional water in the excavation which could pose additional structural concerns.

Granite-Block and/or Concrete Removal

In summary, the anticipated soils to be excavated include granular sand-gravel-silt containing occasional to frequent cobbles and concrete debris (former foundations). Boulders or granite-blocks with larger diameters are possible based on site observations and variation of boring refusals depths.

If large boulders are unable to be removed using large tracked excavators, boulders could be broken up using a hoe ram and then removed using a large excavator. Alternatively, splitting boulders with drilling and chemical splitting methods could be considered. Controlled blasting is not permitted for this project due to the close proximity to existing utilities and foundation structures. Use of a hoe ram or other methods should be performed in conformance to project environmental plans and procedures.

4.2.2 Handling and Interim Storage of PCB Remediation Waste

Interim storage areas will be marked as required by 761.40 and 761.65(c)(9), respectively. Any special requirements for the handling and storage of PCB remediation waste will be determined prior to initiating soil excavation activities. Polyethylene sheeting will be used to protect the ground surface at areas where the material requires temporary storage. The polyethylene sheeting will also be placed over the stockpiled material to serve as a cover system to protect the waste during non-work hours. The storage area will clearly be marked in the field with signs indicating "PCB Remediation Waste". Where practicable, the excavated soil and debris will be directly loaded into waste hauling trucks or roll-off boxes for immediate removal from the site.

4.2.3 Disposal of PCB Remediation Waste – Greater than or equal to 50 ppm

Where practicable (i.e. without requiring stabilization of retaining walls or creating safety concerns for site workers), remediation waste ≥50 ppm will be excavated and removed from the area. All TSCA waste streams with PCB concentrations ≥50 ppm will be transported by licensed hazardous waste haulers and disposed of at Waste Management's Model City facility in Model City, New York (or other similarly licensed facility based upon ability to accept materials at the time of shipment). The appropriate TSCA notification of generation of PCB remediation waste will be filed with EPA, as required.

4.2.4 Disposal of PCB Remediation Waste – Less than 50 ppm

TSCA PCB remediation waste with concentrations less than 50 ppm will be transported by licensed waste haulers to the Juniper Ridge Landfill in Old Town, Maine. Documentation of disposal, in the form of weigh slips, will be maintained with all other documentation for this cleanup.

4.3 Verification Sampling and Analysis

Following the removal of fill and debris from the excavation area, verification samples will be collected. A diagram showing the proposed sampling areas and excavation area is illustrated on Figure 5.

The Excavation Area will be divided into six separate excavation and sampling areas, designated as Areas 1, 2, 3, 4, 5, and 6 as shown. Areas 1, 2, 3, 4, 5, and 6 will be further subdivided to create 20' x 20' sampling areas.

Each discrete 20 X 20 feet area will be sampled independently for cleanup verification purposes.

- Area 1: 786 square feet (>50 ppm) [12 samples from base of the grids and 28 samples from sidewalls]
- Area 2: 1,858 square feet (>50 ppm) [26 samples from base of the grids and 68 samples from sidewalls]
- Area 3: 190 square feet (top 4' >50 ppm) [3 samples from base of the grid (no sidewalls within this grids)]
- Area 4: 6,703 square feet (top 4' <50 ppm) [57 samples from base of the grids and 47 samples from sidewalls]
- Area 5: 385 square feet (top 2' <50 ppm) [4 samples from base of the grids and 9 samples from sidewalls]

Area 6: 340 square feet (removal from 1-9' based on geotechnical recommendations >50 ppm) [4 samples from base of the grids and 2 samples from sidewalls]

The following sample procedures will be conducted:

- Collect three individual samples from the base of each excavation sub-area;
- If a grid has a sidewall, the sidewalls will be sampled in accordance with Subpart O (i.e. 1 sample per 25 feet²);
- Use coordinate-based random sampling to select the sampling locations of the three samples from each excavation sub-area;
- Use a core sampler with a diameter ≥ 2 cm and ≤ 3 cm; and
- Collect samples from the base of the excavation to a maximum depth of 7.5 cm.

All verification samples will be submitted to a certified laboratory for analysis of total PCBs using soxhlet extraction. Quality assurance/quality control (QA/QC) samples will be collected at a rate of one per twenty samples.

To demonstrate compliance with the target cleanup level of 25 ppm, all individual sample results from the verification samples collected from the base/sidewall of the excavation must be equal to or less than 25 ppm.

If specific sub areas within the excavation zones indicate PCB levels greater than the target cleanup level of 25 ppm, we are prepared to complete additional soil excavation and disposal and verification sampling if practicable (i.e., if there are no structural concerns related to the railroad or retaining walls, groundwater entering the excavation and creating a potentially hazardous situation for site workers).

4.4 Backfill and Site Restoration

Following receipt of acceptable (≤ 25 ppm total PCB) post-excavation verification sample results, the excavations will be backfilled and restored with an In-Situ Cap System. This cap system is to be constructed over remaining PCB impacted material to prevent human exposure, manage infiltration of water, and minimize erosion of soil.

Prior to construction of cap system, a borrow source characterization will be conducted to document that proposed barrier soil meets the requirements set forth in 40 CFR §761.75(b)(1)(ii) through (b)(1)(v), and to determine acceptable construction methods capable of attaining in-place permeability requirements. Understanding that PCB concentrations remaining at the Site may be up to 25 ppm, an enhanced capping system is being proposed that is significantly more protective than required by EPA's PCB remediation rules.

Twelve inches of barrier soil will be placed from 3-4 feet below grade over remaining impacted soil. A 40-mil textured high density polyethylene (HDPE) geomembrane liner will overly the barrier soil layer and be installed in accordance with manufacturers recommendations by an experienced installer. A drainage geocomposite layer comprised of bi-planar HDPE netting sandwiched between non-woven geotextile filter fabric will overlay the geomembrane. Geocomposite will extend over the limit of the cap and be installed according to manufacturer's recommendations by an experienced installer. In addition to transmitting surface water over the cap system, the HDPE netting and geotextile fabric provides a protective layer over the geombrane.

A 12 inch thick compacted sand drainage layer shall be placed over the installed geocomposite. A crushed stone underdrain will be installed at the eastern limit of the cap system and extend along the existing granite retaining wall adjacent to the Saco River to allow surface water to drain off of the cap system.

After the installation of drainage sand layer and stone underdrain, a marker layer (orange site safety fence) is to be installed at a depth of 2 feet below the ground surface to alert persons of the presence of underlying materials that should not be disturbed. An 18 inch thick layer of granular borrow is to be placed and compacted over the marker layer.

The surface of compacted granular borrow is to be scarified to a depth of 2 inches prior to placement of 6 inches of loam then seeded, and mulched.

4.5 Post-Construction Activities

A post-construction report will be prepared following remediation to document the completed field activities, present the verification sampling data, and provide copies of the waste manifests executed for the transportation and disposal of the waste. All reports will be maintained on file at the owner's property and at City Hall, in accordance with the record-keeping requirements of Subpart J of 40 CFR 761. Copies of such records will be made available to the EPA, upon request.

Land use in the cleanup area and an abutting 25 feet buffer will be subject to restrictions contained in a deed notice and environmental covenant. Execution of the environmental covenant will be completed within 60-days of the agencies' "certificate of completion" or equivalent. Recordation of the Deed Covenant as well as other Institutional controls (non-PCB related) will be documented in the MEDEP VRAP.

5.0 RISK ANALYSIS

Currently, the area of concern is undeveloped. The ground surface consists of asphalt paving and grass/meadow. Due to the industrial use of the property since construction of the waste to energy facility, public access to the area of concern has been prohibited. As a result, exposure scenarios are limited and individuals would spend only limited time in the area, well below the hourly threshold for "low occupancy" use.

However, the property has been purchased by the City of Biddeford, the facility is being decommissioned and future plans for the area of concern call for a park setting. Given this change of use, the current distribution of PCB impacts will pose an unacceptable exposure risk.

PCB concentrations in the 0-2 feet below grade interval range up to 631 ppm. The 0-2 feet below grade interval is considered a Direct Contact interval where workers involved in constructing the park, as well as future park users could come into direct contact with PCB impacted fill. The park setting is considered to be a "high occupancy" setting and exposure to PCB concentrations above 1 ppm are considered to pose an unacceptable risk.

In addition, there are no restrictions for use of the property and land use changes could result in further disturbance, redistribution of impacted fill and exposure scenarios that would potentially increase risk to construction workers, future occupants and the environment.

The proposed Remediation Plan includes a tiered risk reduction approach that is protective of human health and the environment and includes the following risk reduction measures:

- 1) Removal of fill materials containing greater than 25 ppm from the area of concern to a depth of 4 feet, well below the Direct contact interval.
- 2) Installing an enhanced capping system that extends from 4 feet below grade to ground surface. Cap materials include barrier soils, geomembrane liners, geocomposite and drainage layers, a marker layer and 2 feet of clean fill and loam to prevent exposure to PCB impacted fill.
- 3) Execution of a Deed Covenant to restrict land use and prohibit activities that could result in disturbance of fill below the cap system. Prior to any ground disturbance within the limits of the cap, the City Environmental Office will be contacted to obtain authorization.
- 4) Development of a long-term maintenance program to inspect the area, maintain the integrity of the capping system and provide awareness training to City personnel responsible for activities in the park.

In combination, these measures will result in an elimination of exposure pathways through physical barriers and Institutional Controls. Long-term maintenance plans will include an awareness and communication program to reinforce the restrictions and ensure protection of human health and the environment.

6.0 REFERENCES

- Credere Associates, LLC (Credere). September 18, 2012. "Phase I Environmental Site Assessment, Maine Energy Recovery Company Property, 3-11 Lincoln Street, Biddeford, Maine."
- Credere. October 25, 2012. "Phase II Environmental Site Assessment, Maine Energy Recovery Company Property, 3-11 Lincoln Street, Biddeford, Maine."
- Credere. November 15, 2012. "Supplemental Phase II Environmental Site Assessment, Maine Energy Recovery Company Property, 3-11 Lincoln Street, Biddeford, Maine."
- Credere. November 5, 2012. "Draft Voluntary Response Action Program Work Plan."
- EDR. 2012. "Sanborn Fire Insurance Maps, 1867-1976 Maine" Sanborn. 1932, 1947, 1976.
- EDR. 2012. "1956, 1960, 1970, 1975, 1986, 1990, 1998, 2003 Aerial Photographs.

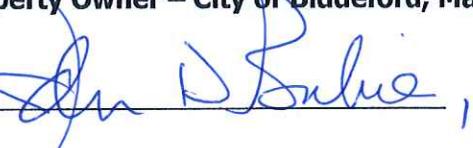
CERTIFICATION

The undersigned owner of the property where the cleanup site is located and the party conducting the cleanup certify that all sampling plans, sampling collection procedures, extraction procedures and instrumental/chemical analysis procedures used to assess or characterize the PCB contamination at the cleanup site are on file at the location indicated below and are available for EPA inspection, as set forth below:

Document Location

City of Biddeford
205 Main Street
Biddeford, Maine 04005

Property Owner – City of Biddeford, Maine


John D. Bubier,
Authorized Signature

4/16/2014
Date

John D. Bubier
Name of Authorized representative (print)

City Manager
Title

Party Conducting the Cleanup – Maine Energy


Brian Oliver
Authorized Signature

4/16/14
Date

Brian Oliver
Name of Authorized representative (print)

Vice President
Title

Maine Energy PCB Investigation - Table 1
October 2012 Phase II ESA

LOCATION	CA-SB-3	CA-SB-5	CA-SB-6	CA-SB-6	CA-SB-7	CA-SB-8	CA-SB-8	CA-SB-9	CA-SB-10	CA-SB-11	CA-SB-12	CA-SB-13	TSCA
DATE	10/2/12	10/3/12	10/3/12	10/3/12	10/2/12	10/3/12	10/2/12	10/3/12	10/3/12	10/3/12	10/2/12	10/2/12	
DEPTH	10-12.5'	5-10'	0-5'	12.5-15'	7.5-10'	0-2.5'	5-10'	0-4'	10-15'	10-15'	15-18'	10-12'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	ND	250	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
TOTAL PCBs:	ND	250	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1

LOCATION	CA-SS-1	CA-SS-2	CA-SS-3	CA-SS-4	CA-SS-5	TSCA
DATE	10/3/12	10/3/12	10/3/12	10/3/12	10/3/12	
DEPTH	0-6"	0-6"	0-6"	0-6"	0-6"	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	
PCB - 1260	ND	ND	ND	0.750	ND	
TOTAL PCBs:	ND	ND	ND	0.750	ND	1

Table Notes:

ND = Value is below method detection limits

mg/kg = milligrams per kilogram, or parts per million

TSCA = USEPA Toxic Substance Control Act

Table 2
PCB SAMPLE RESULTS
Maine Energy Recovery Company Property
3-11 Lincoln Street, Biddeford, Maine

Lab Sample #	Boring # (CA-SB-xx)	Sample Depth	Method 4020 Semiquantitative Result							4020 batch #	QA Samples soil spikes at x ppm		
			<1	~1	1 to 10	~10	10 to 50	~50	>50		0	1	10
1	5R	0-6"			x					1	<1	1-10	1-10
2	5R	6"-2.5'					x	x		1			
3	5R	2.5-5'					x	x		1			
4	105	0-6"				r	x			1			
5	105	6"-3'	x r							1			
6	111	0-6"			x					1			
7	111	6"-2.5'					x r			1			
8	111	2.5-5'					x r			1			
9	111	5-7.5'						IV		1			
10	111	7.5-10'	x							1			
11	110	0-6"			x			x r		1			
12	110	5-7.5'								1			
13	110	7.5-10'	x							1			
14	113	0-6"			x					2	<1	~1	1-10
15	113	6"-2.5'			x					2			
16	113	2.5-5'			r	x				2			
17	113	5-7.5'			x					2			
18	113	7.5-10'	x							2			
19	103	0-6"			x		x			2			
20	103	6"-2.5'			x		x			2			
21	103	2.5-5'			x					2			
22	103	5-7.5'	x							2			
23	103	7.5-10'	x							2			
24	114	0-6"				r	x			2			
25	114	6"-2.5'	x							3	<1	1-10	1-10
26	114	2.5-5'	x r			x				3			
27	114	5-7.5'			x					3			
28	114	7.5-10'	x							3			
29	106	0-6"		x					xx	3			
30	106	6"-2.5'		x						3			
31	106	2.5-5'		x		x				3			
32	106	5-6'		x						3			
33	112	0-6"			x					3			
34	112	6"-2'			IV					3			
35	100	0-6"			x					4	<1	1-10	1-10
36	100	6"-2.5'			x		x			4			
37	100	2.5-5'			x					4			
38	100	5-7.5'			x			xx		4			
39	100	7.5-10'			x					4			
40	101	0-6"							x	4			
41	101	6"-5'			x				x	4			
42	101	5-10'			xx					4			
43	102	0-6"			x					4			
44	102	6"-2.5'			x					4			
45	102	2.5-5'			x				x r	4			
46	104	0-6"			xx					4			
47	104	5-7.5'			x					5	<1	1-10	1-10
48	104	7.5-10'			x					5			
49	107	0-6"		x				r	xx	5			
50	107	6"-2.5'		x						5			
51	107	2.5-5'		x		x				5			
52	107	5-7.5'		x						5			
53	107	7.5-10'		xx						5			
54	108	0-6"					xx			5			
55	108	6"-2.5'			x					5			
56	108	2.5-3'			xx					5			
57	109	0-6"			x					6			
58	109	6"-2.5'			x					6			
59	109	2.5-5'		xx		x				6			
60	109	5-7.5'		x						6			
61	109	7.5-10'		x						6			
62	115	0-6"		xx			x			6			
63	115	6"-2.5'								6			
64	116	0-6"		xx						6			
65	116	6"-3'								6			
66	116	2.5-5'							x	6			
67	116	5-7.5'					x			6			
68	116	7.5-10'	x		x					6			
69	117	0-6"			x					6			

Notes

This results table was provided by CAPE Technologies and the data has not been altered

Analysis completed by CAPE Technologies on November 9, 2012

All reported results on this table were analyzed via EPA Method 4020

Each batch included a zero plus calibrators at 1, 10, and 50 ppm

Results are given as the semiquantitative result relative to the calibrators in that batch

XX - Replicates (parallel subsample in same batch) are shown as two "x" marks on one sample line

R - Repeat analyses (later analysis of 2nd aliquot of extract) are shown as an "r" on the same line as the original ("x")

QA samples consisting of soils spiked as noted were extracted and an aliquot of each extract was analyzed in each batch

IV - Samples 8 & 34 (ital. & underlined) were cloudy due to high levels of coextracted hydrocarbons and results should be considered invalid

Maine Energy PCB Investigation - Table 3

October 2012 Supplemental Phase II ESA

LOCATION	CA-SB-110	CA-SB-111	CA-SB-111	CA-SB-111	TSCA
DATE	10/25/12	10/25/12	10/25/12	10/25/12	
DEPTH	7.5-10'	2.5-5'	5-7.5'	7.5-10'	
	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	1
PCB - 1221	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	
PCB - 1260	ND	6300	55	0.05	
TOTAL PCBs:	ND	6300	55	0.05	1

Table Notes:

ND = Value is below method detection limits

mg/kg = milligrams per kilogram, or parts per million

TSCA = USEPA Toxic Substance Control Act

Maine Energy PCB Investigation - Table 4
 January 2013 Results

LOCATION	SE-SB-206	SE-SB-206	SE-SB-207	SE-SB-207	SE-B-207	SE-SB-208	SE-SB-208	SE-SB-208	SE-SB-208	SE-SB-209	SE-SB-209	SE-SB-209	TSCA
DATE	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	
DEPTH	3"-2'	2-12'	3"-3'	3-4.5'	5-6'	3"-4'	4-6'	6-8'	8-10"	2"-3'	3-9'	9-12.5'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	1.7	0.31	ND	6.3	2.30	ND	2.5	5.0	25.0	ND	3500	0.13	
TOTAL PCBs:	1.7	0.31	ND	6.3	2.30	ND	2.5	5.0	25.0	ND	3500	0.13	1

LOCATION	SE-SB-212	SE-SB-201	SE-SB-201	SE-SB-202	SE-SB-202	SE-SB-203	SE-SB-203	SE-SB-203	SE-SB-204	SE-SB-204	SE-SB-205	TSCA
DATE	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	1/10/2013	
DEPTH	2-20"	3"-4'	4-5"	4"-2.5'	4-5"	2"-2.5'	4-5'	5-7.5'	3"-3"	3-5"	3"-3'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.194	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	1.70	ND	0.40	ND							
PCB - 1260	7.1	ND	4.60	ND	ND	0.05	85	11	ND	11	0.07	
TOTAL PCBs:	7.1	1.700	4.60	ND	ND	0.05	85	11	ND	11	0.07	1

LOCATION	SE-SB-205	SE-SB-205	SE-SB-210*	SE-SB-211	SE-SB-211	TSCA
DATE	1/10/13	1/10/13	1/10/13	1/10/13	1/10/13	
DEPTH	3.5-7.5'	10.5-12.5'	0-20"	2"-3.5'	4-9"	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	1
PCB - 1248	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	
PCB - 1260	0.3	ND	ND	0.230	48	
TOTAL PCBs:	0.3	ND	ND	0.230	48	1

Table Notes:

ND = Value is below method detection limits

mg/kg = milligrams per kilogram, or parts per million

TSCA = USEPA Toxic Substance Control Act

Maine Energy PCB Investigation - Table 5
February 2013 Results

LOCATION	SE-SB-213	SE-SB-213	SE-SB-213	SE-SB-213	SE-SB-214	SE-SB-214	SE-SB-214	SE-SB-214	SE-SB-215	SE-SB-215	SE-SB-215	SE-SB-215	SE-SB-215	TSCA
DATE	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	
DEPTH	0-2'	2-4'	4-6'	6-7.5'*	0-2'	2-4'	4-6'	6-7'*	0-2'	2-4'	4-6'	6-8'	8-10*	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	0.707	0.585	1.79	9.14	0.385	1.280	17.20	8.16	ND	3.21	2.26	3.22	44.6	
TOTAL PCBs:	0.707	0.585	1.79	9.14	0.385	1.280	17.20	8.16	ND	3.21	2.26	3.22	44.6	1

LOCATION	SE-SB-216	SE-SB-216	SE-SB-216	SE-SB-217	SE-SB-217	SE-SB-217	SE-SB-217	SE-SB-219	SE-SB-219	SE-SB-219	SE-SB-219	SE-SB-219	SE-SB-219	TSCA
DATE	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	
DEPTH	2.5"-24"	2-4'	4-6*	0-2'	2-4'	6-8'	8-10"-	0-2'	2-4'	4-6'	6-8'	8-10*		
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	0.065	12.90	ND	ND	1.160	0.853	6.68	ND	0.380	11.0	39.8	54.1		
TOTAL PCBs:	0.65	12.90	ND	ND	1.160	0.853	14.54	ND	0.380	11.0	39.8	54.1		1

LOCATION	SE-SB-220	SE-SB-220	SE-SB-220	SE-SB-220	SE-SB-220	SE-SB-220	SE-SB-221	SE-SB-221	SE-SB-221	SE-SB-224	SE-SB-224	SE-SB-224	SE-SB-224	TSCA
DATE	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	2/19/13	
DEPTH	0-2'	2-4'	4-6'	6-8'	8-10'	10-11*	0-2'	2-4'	4-5*	0-2'	2-4'	6-8*		
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	ND	0.284	2.61	13.90	19.80	489.0	0.047	0.308	0.068	9.55	0.229	9.26		
TOTAL PCBs:	ND	0.284	2.61	13.90	19.80	489.0	0.047	0.308	0.068	9.55	0.229	9.26		1

LOCATION	SE-SB-225	SE-SB-225	SE-SB-225	SE-SB-225	TSCA
DATE	2/19/13	2/19/13	2/19/13	2/19/13	
DEPTH	0-2'	2-4'	4-6'	6-7.5*	
	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	
PCB - 1260	0.31	15.8	17.6	10.90	
TOTAL PCBs:	0.31	15.8	17.6	10.90	1

Maine Energy PCB Investigation - Table 5

February 2013 Results

LOCATION	SE-SB-218	SE-SB-218	SE-SB-222	SE-SB-222	SE-SB-222	SE-SB-222	SE-SB-223	SE-SB-223	SE-SB-223	SE-SB-223	SE-SB-223	SE-SB-226	SE-SB-226	SE-SB-226	TSCA
DATE	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	
DEPTH	0'-2'	2'-4'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-5"
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	0.328	0.696	2.74	0.447	0.823	0.146	0.237	0.133	179	3.92	3.93	8.00	631	9660	3650
TOTAL PCBs:	0.328	0.696	2.74	0.447	0.823	0.146	0.237	0.133	179	3.92	3.93	8.00	631	9660	3650

LOCATION	SE-SB-227	SE-SB-227	SE-SB-227	SE-SB-227	SE-SB-228	SE-SB-228	SE-SB-228	SE-SB-228	SE-SB-229	SE-SB-229	SE-SB-229	SE-SB-229	SE-SB-229	SE-SB-230	SE-SB-230	TSCA	
DATE	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13		
DEPTH	0'-2'	2'-4'	4'-6'	6'-8'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		
PCB - 1260	2.84	1.93	0.996	4.32	86.2	1.36	19.5	3.98	15.50	577	192	50.6	149	26.3	27.7	3.08	
TOTAL PCBs:	2.84	1.93	0.996	4.32	86.2	1.36	19.5	3.98	15.50	577	192	50.6	149	26.3	27.7	3.08	

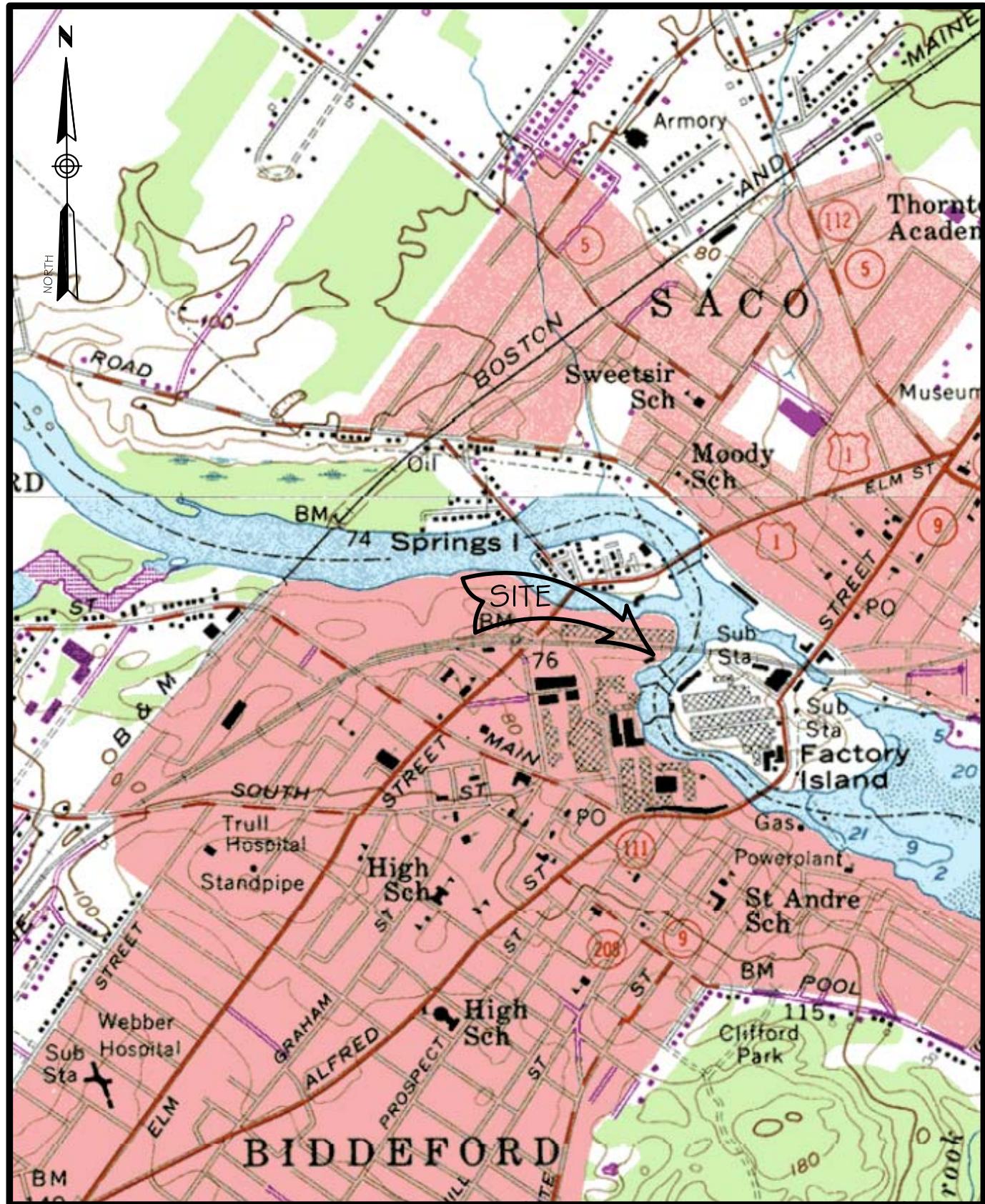
LOCATION	SE-SB-231	SE-SB-231	SE-SB-231	SE-SB-231	SE-SB-231	SE-SB-232	SE-SB-232A	SE-SB-232A	SE-SB-232A	SE-SB-232A	SE-SB-233	SE-SB-233	SE-SB-233	SE-SB-233	SE-SB-233	TSCA
DATE	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	
DEPTH	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	17.8	324	164	3150	2330	553	4.49	4.34	5.09	0.241	6.30	1.02	0.267	0.404	1	
TOTAL PCBs:	17.8	324	164	3150	2330	553	4.49	4.34	5.09	0.241	6.30	1.02	0.267	0.404	1	

LOCATION	SE-SB-234	SE-SB-234	SE-SB-234	SE-SB-234	SE-SB-235	SE-SB-235	SE-SB-235	SE-SB-235	SE-SB-236	SE-SB-236	SE-SB-236	SE-SB-236	SE-SB-236	SE-SB-236	SE-SB-236	TSCA
DATE	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	2/20/13	
DEPTH	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	0'-2'	2'-4'	4'-6'	6'-8'	8'-10'	
	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	mg/kg	
PCB - 1016	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1221	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1232	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1242	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1248	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1254	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
PCB - 1260	25.10	2.35	1.84	2.26	1.54	2.24	0.488	ND	1.38	0.399	0.174	0.67	2.24	2.12	0.975	0.735
TOTAL PCBs:	25.10	2.35	1.84	2.26	1.54	2.24	0.488	ND	1.38	0.399	0.174	0.67	2.24	2.12	0.975	0.735

Table Notes:

ND = Value is below method detection limits
mg/kg = milligrams per kilogram, or parts per million
TSCA = USEPA Toxic Substance Control Act
* = Denotes refusal at depth
- = Hole collapsed upon removal of tooling at 10'

Figures



SITE LOCATION PLAN

MAINE ENERGY RECOVERY COMPANY
PCB INVESTIGATION

DATE: 3/26/2013

DRAWN BY: TND

CHECKED BY: JKC

JOB: 12-3259.1

NOT TO SCALE

FILE: 3259.1 FIG.dwg



FIGURE

640 MAIN ST.
LEWISTON, ME 04240

Tel.: (207) 795-6009
Fax: (207) 795-6128

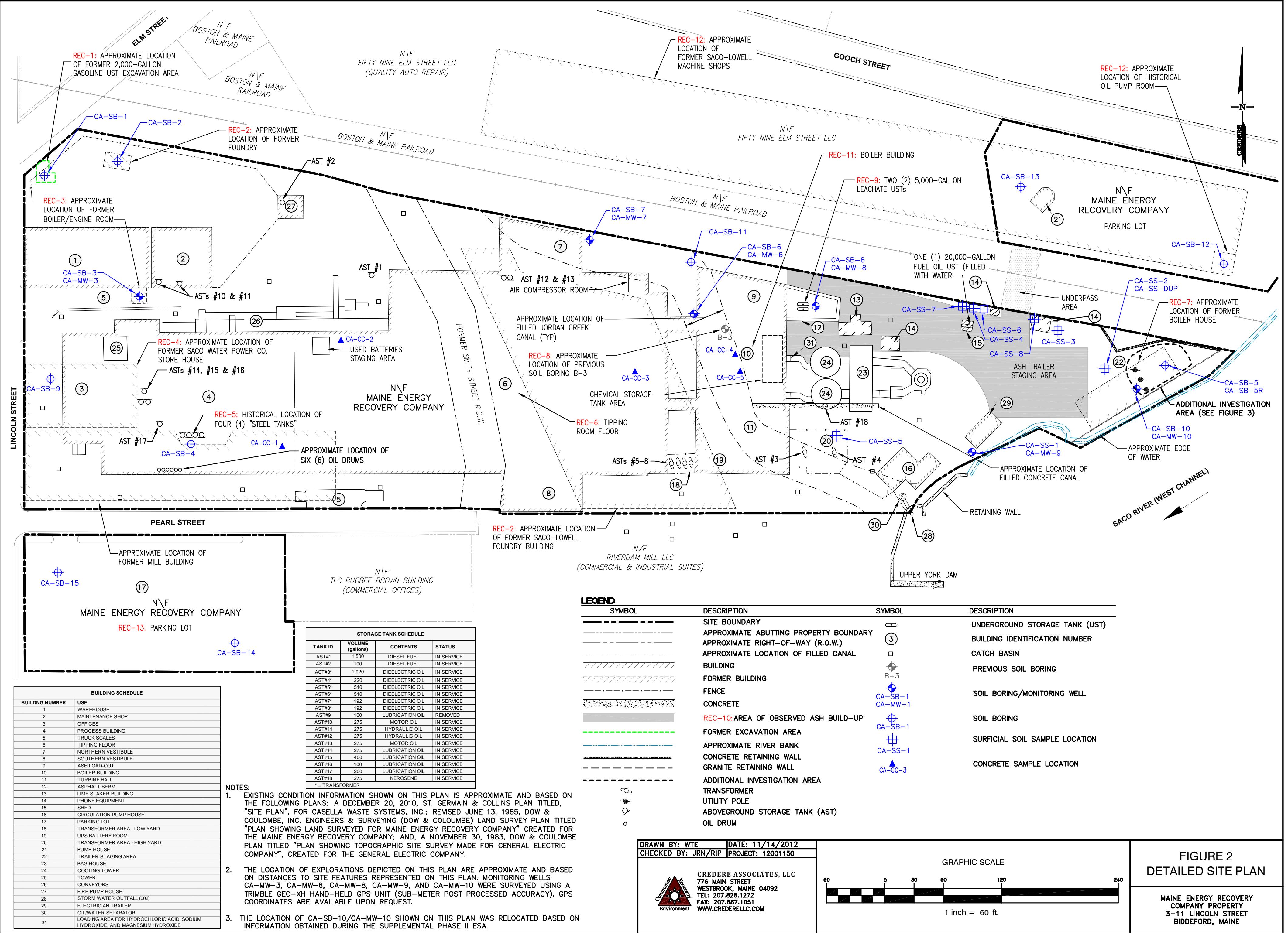
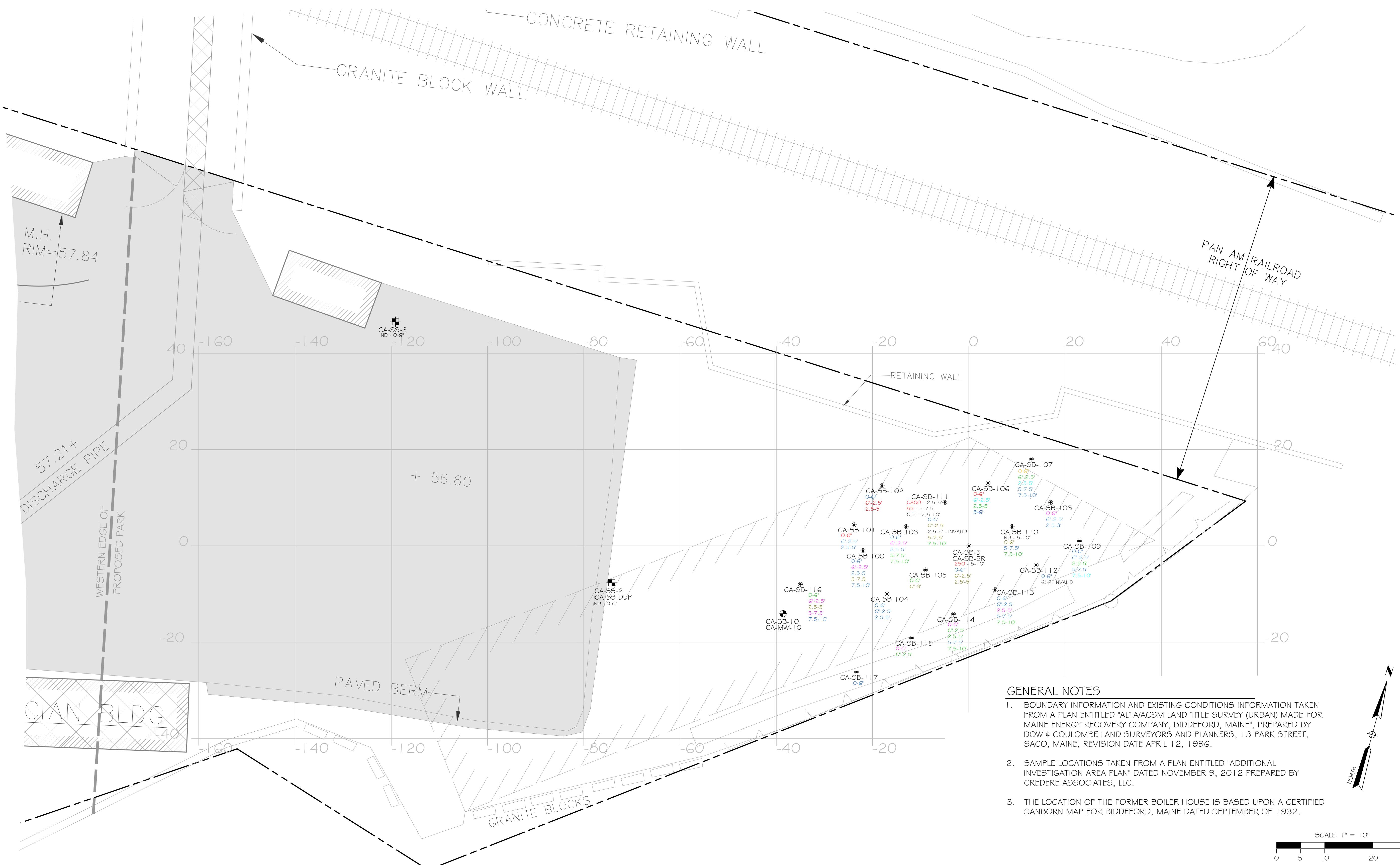


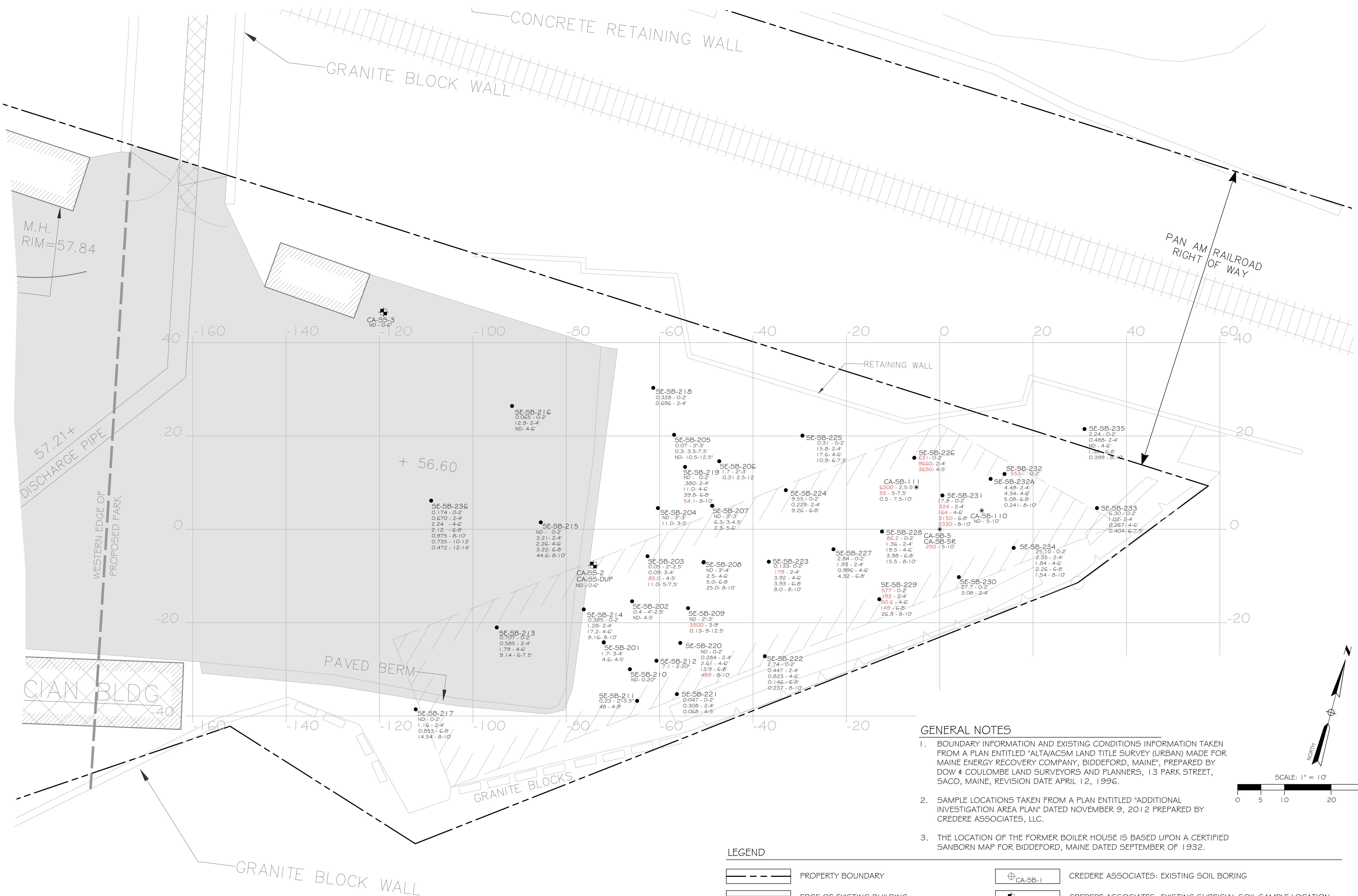
FIGURE 2 DETAILED SITE PLAN



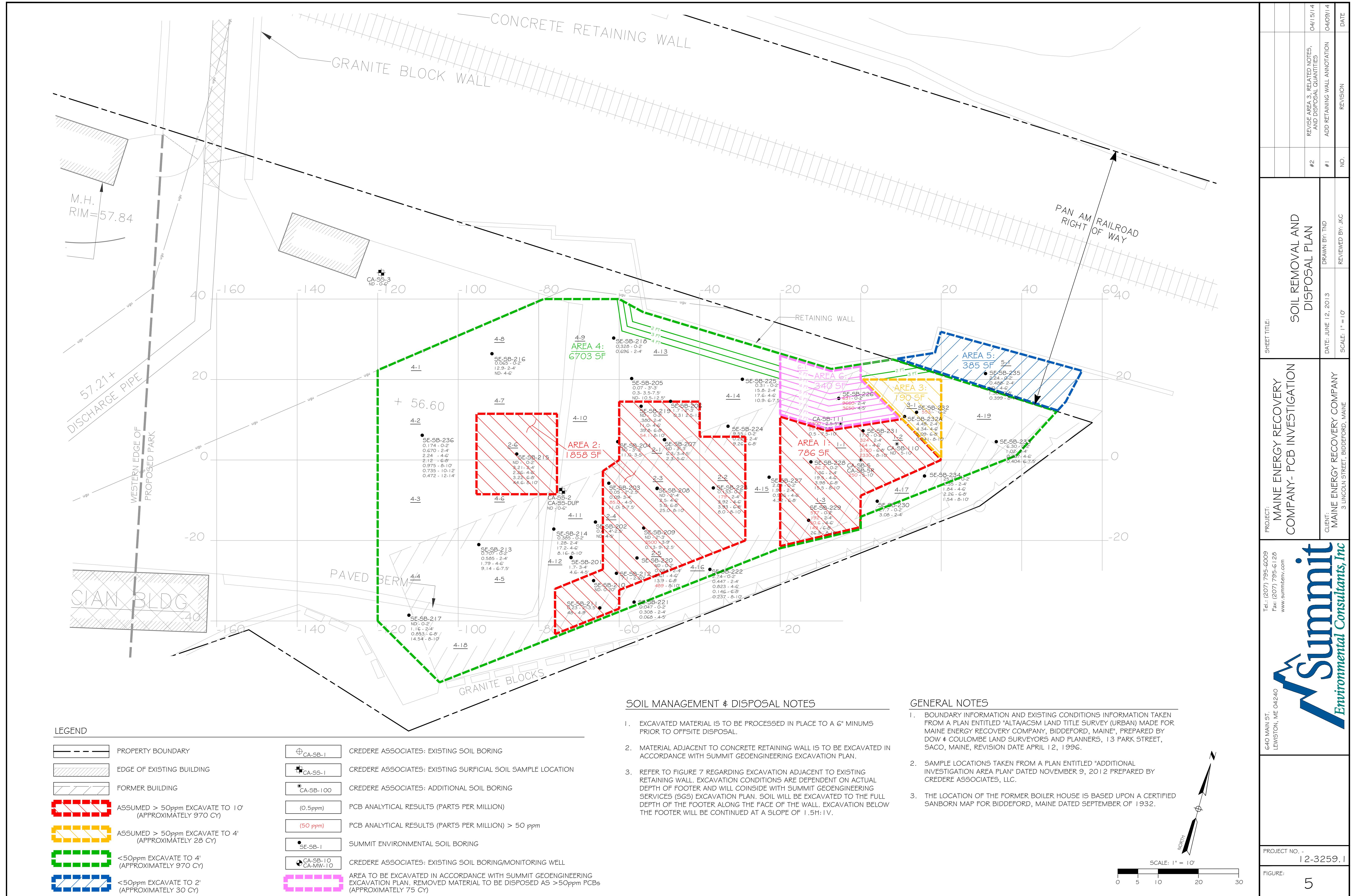
LEGEND

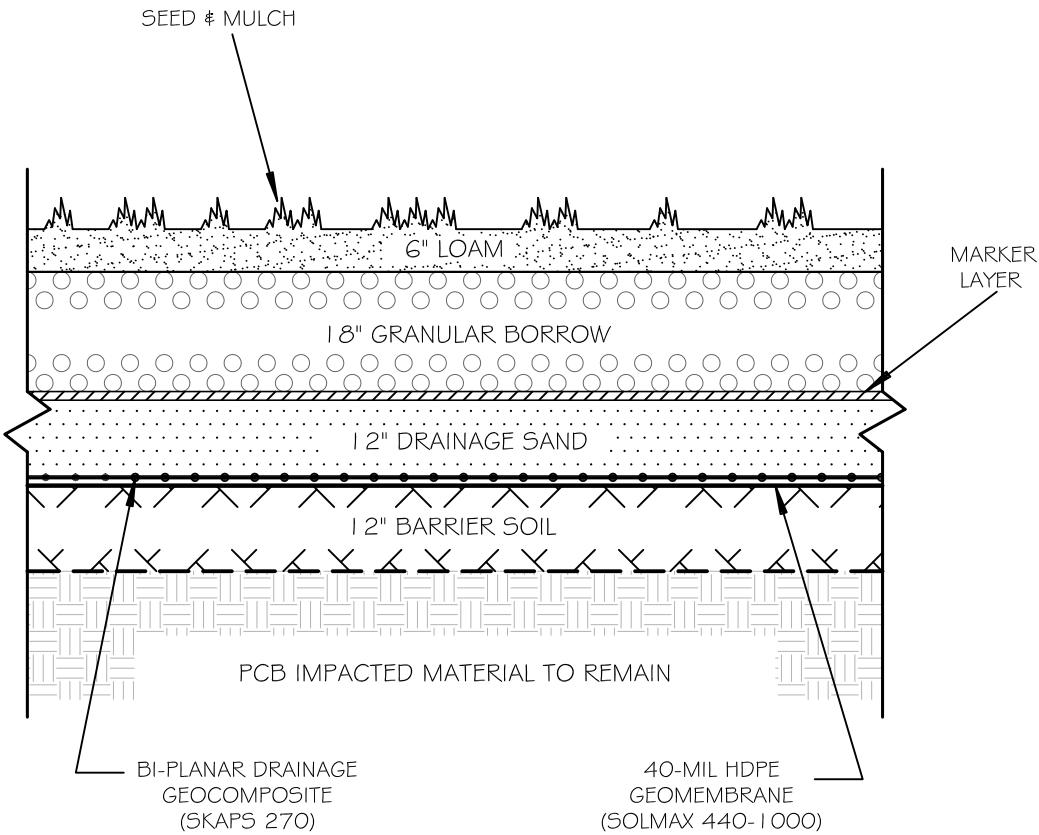
PROPERTY BOUNDARY	CA-55-1	CREDERE ASSOCIATES: EXISTING SURFICIAL SOIL SAMPLE LOCATION	O-G*	<1 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
EDGE OF EXISTING BUILDING	CA-SB-100	CREDERE ASSOCIATES: ADDITIONAL SOIL BORING	O-G*	~1 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
FORMER BUILDING	(0.5ppm)	PCB ANALYTICAL RESULTS (PARTS PER MILLION)	O-G*	1 TO 10 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
PCB INVESTIGATION AREA OF CONCERN	(50 ppm)	PCB ANALYTICAL RESULTS (PARTS PER MILLION) > 50 ppm	O-G*	~10 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
ASPHALT STAGING AREA	CA-SB-10 CA-MW-10	CREDERE ASSOCIATES: EXISTING SOIL BORING/MONITORING WELL	O-G*	10 TO 50 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
			O-G*	~50 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT
			O-G*	>50 ppm PCBs AS DETERMINED BY CAPE TECHNOLOGIES IMMUNOASSAY SCREENING KIT

PROJECT NO. -	12-3259.1	FIGURE:	3
640 MAIN ST. LEWISTON, ME 04240	Tel.: (207) 795-6009 Fax: (207) 795-6128 www.summitenv.com	PROJECT: MAINE ENERGY RECOVERY COMPANY- PCB INVESTIGATION	SHEET TITLE: SAMPLE LOCUS PLAN: CREDERE ASSOCIATES
CLIENT: MAINE ENERGY RECOVERY COMPANY 3 LINCOLN STREET, BIDDEFORD, MAINE	DATE: JANUARY 11, 2013 SCALE: 1" = 10'	DRAWN BY: TND	REVIEWED BY: JJC
02	RELOCATE BOILER HOUSE/SANBORN	01	ADD. BORINGS/MTS SURVEY DATA
03/04/13	3/22/13	NO.	REVISION
DATE			



SAMPLE LOCUS PLAN		SAMPLE LOCUS PLAN	
PROJECT:	MAINE ENERGY RECOVERY COMPANY- PCB INVESTIGATION	DRAWN BY: TND	REVIEWED BY: JJC
CLIENT:	MAINE ENERGY RECOVERY COMPANY 3 LINCOLN STREET, BIDDEFORD, MAINE	DATE: JANUARY 11, 2013	SCALE: 1" = 10'
640 MAIN ST. LEVINGTON, ME 04240	Tel: (207) 795-6009 Fax: (207) 795-6128 www.summitenv.com	O2 RELOCATE BOILER HOUSE/SANBORN	3/04/13
		O1 ADD. BORINGS/MTS SURVEY DATA	2/22/13
		NO.	REVISION
			DATE





NOTES:

1. BARRIER SOIL:
 - HYDRAULIC CONDUCTIVITY: $\leq 1 \times 10^{-7}$ cm/s²
 - GRADATION: GREATER THAN 30% PASSING THE #200 SIEVE
 - LIQUID LIMIT: ≥ 30
 - PLASTICITY INDEX: ≥ 15
2. GRADE BARRIER SOIL FOR POSITIVE DRAINAGE TOWARDS SACO RIVER BANK AND PROPOSED STONE UNDERDRAIN.

IN-SITU CAP SYSTEM
TYPICAL DETAIL

MAINE ENERGY RECOVERY COMPANY
PCB INVESTIGATION

DATE: 3/26/2013

DRAWN BY: TND

CHECKED BY: JKC

JOB: 12-3259.1

NOT TO SCALE

FILE:3259.1FIG.dwg

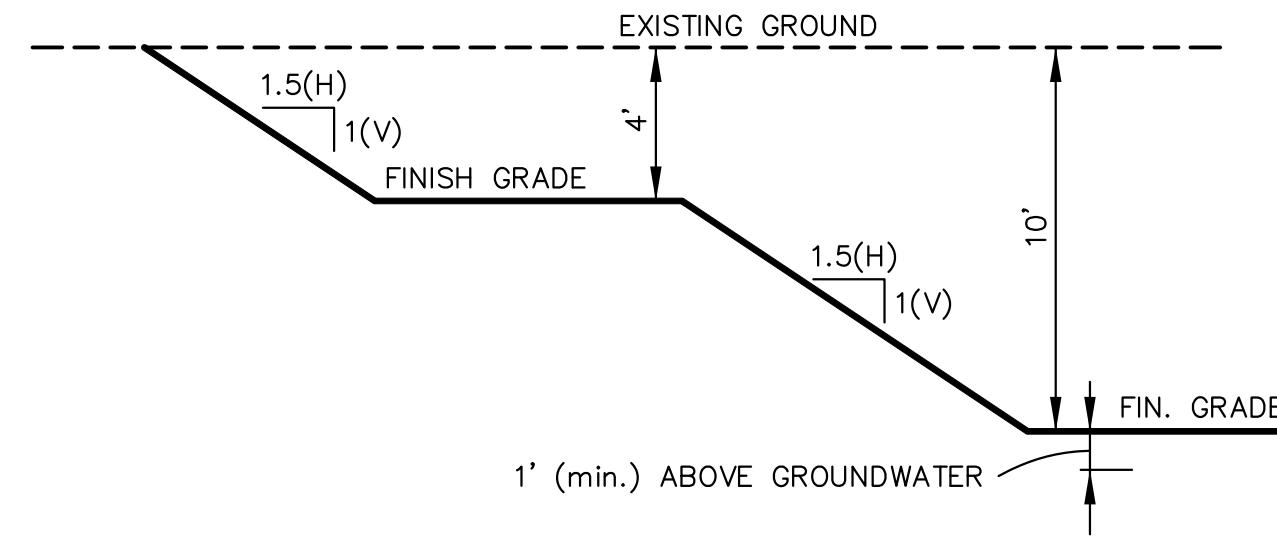
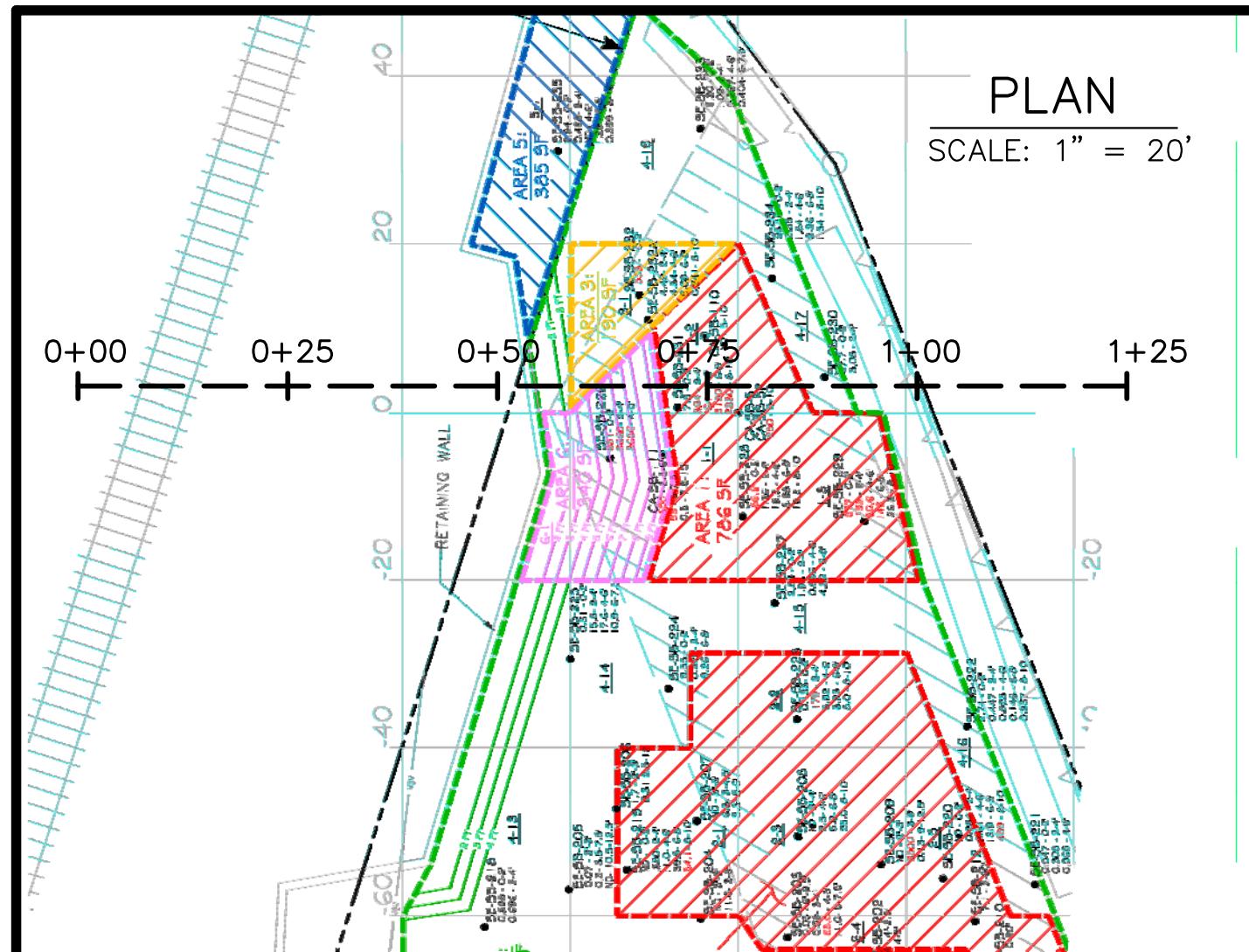


FIGURE

6

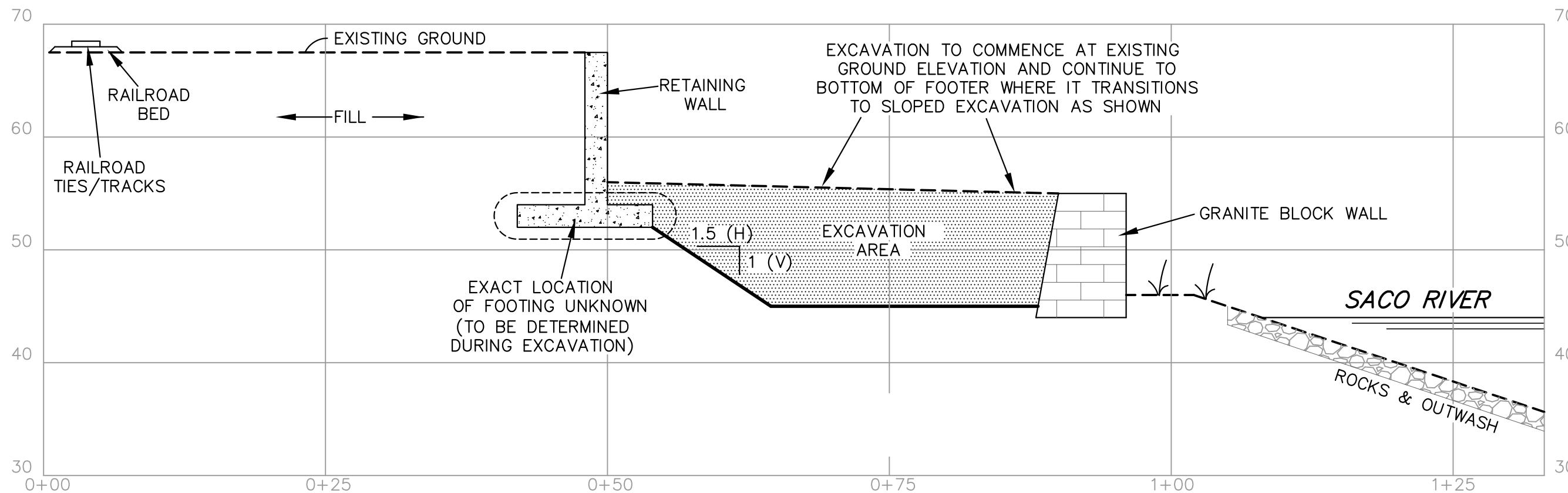
640 MAIN ST.
LEWISTON, ME 04240

Tel.: (207) 795-6009
Fax: (207) 795-6128



EXCAVATION TRANSITION BETWEEN DIFFERING DEPTHS

SCALE: 1" = 5'

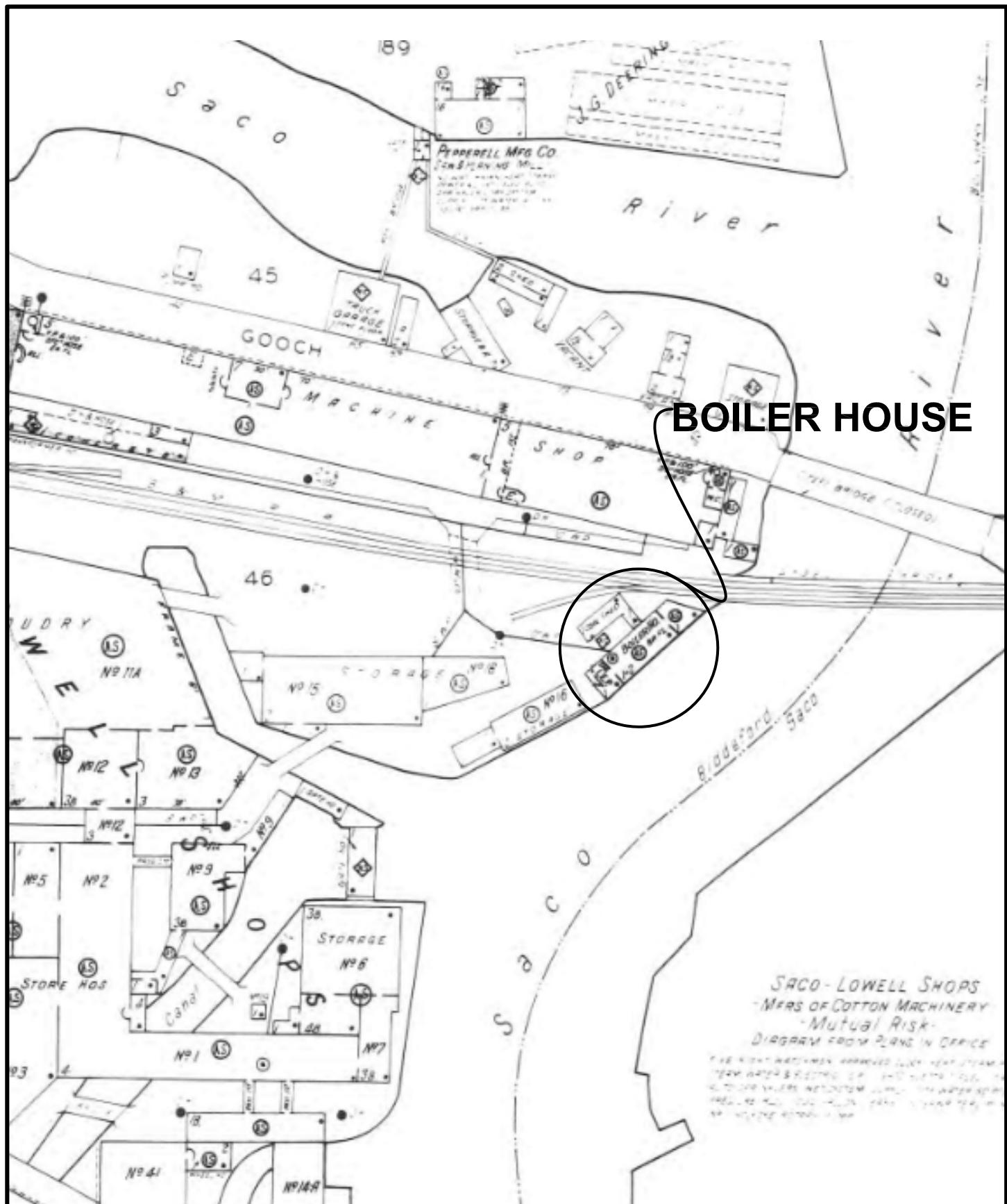


434 CONY ROAD AUGUSTA, ME 04330	Tel.: (207) 318-7761 Fax: (207) 629-9094 www.summitgeoeng.com	FIGURE: 7
PROJ.#: 13051		
4-19-14 - NEW PLAN VIEW REVISED: 4-12-13 - NEW PLAN VIEW BASE MAP		
PROJECT: MAINE ENERGY RECOVERY FACILITY BIDDEFORD, MAINE	CLIENT: SUMMIT ENVIRONMENTAL CONSULTANTS	
DATE: MARCH 29, 2013	APPR. BY: CWC	
AS NOTED	DRAWN BY: KRF	

SUMMIT
GEOENGINEERING SERVICES

Appendix A

Sanborn Fire Insurance Maps and Aerial Photos



PROJECT TITLE:

PCB REMEDIATION BIDDEFORD, MAINE

DWG

BY: RMD

DATE: 1/2014

- REV:

SHEET TITLE:

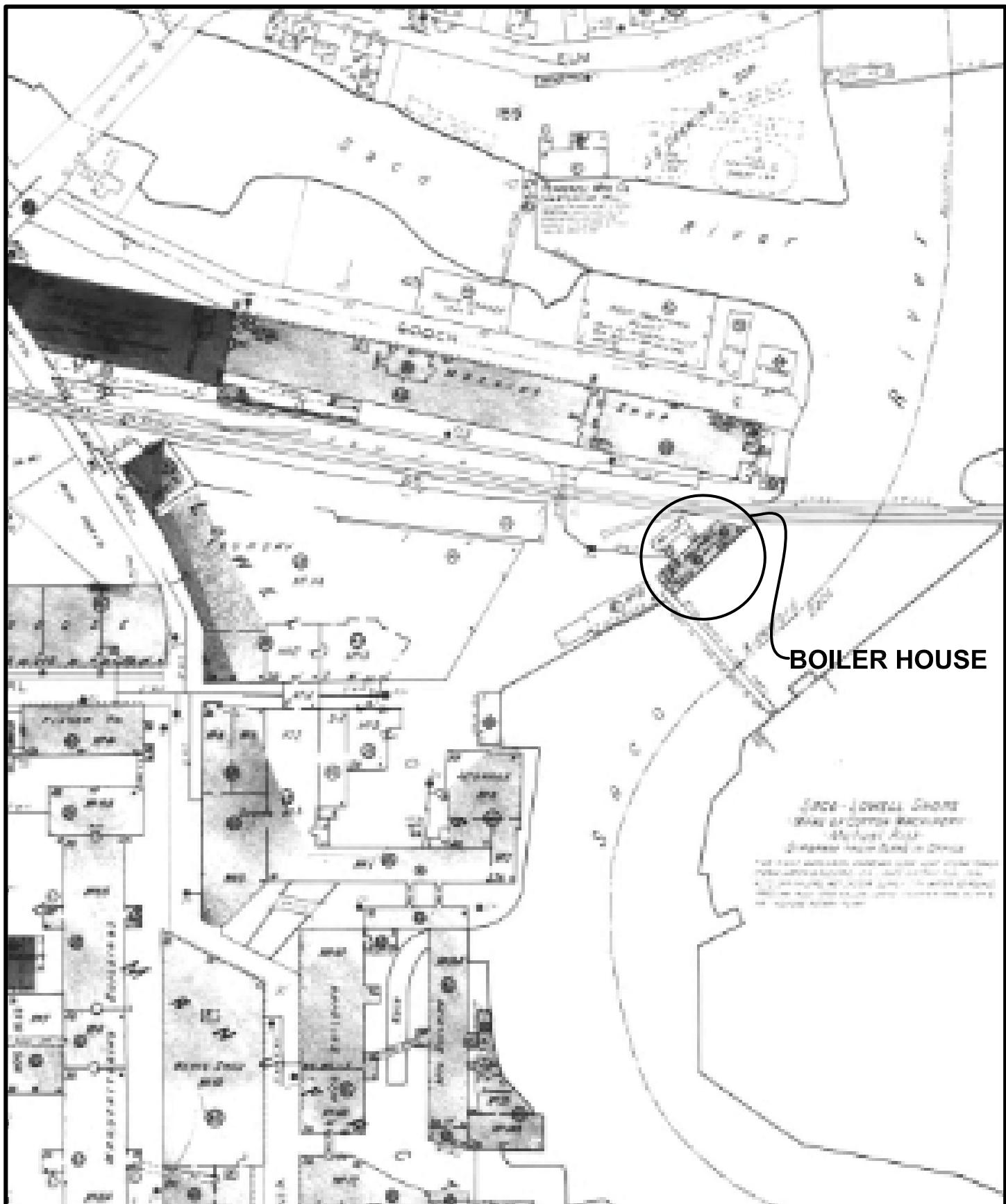
MAINE ENERGY
1932 SANBORN MAP

JN: 10-2240-1 REV: 116

SCALE:

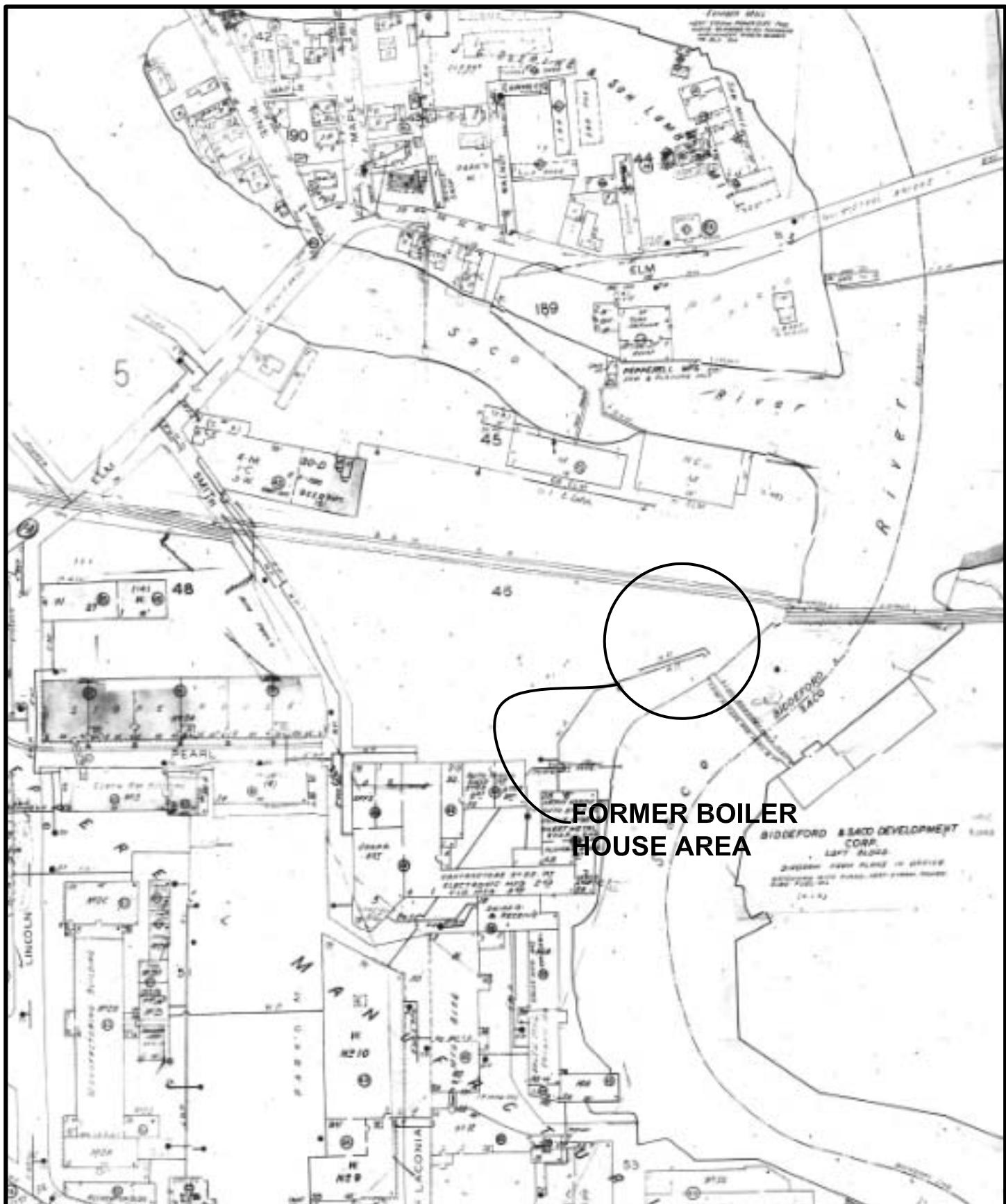
NIS REV DATE: 1/2014





PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A2	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1947 SANBORN MAP	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





PROJECT TITLE:

**PCB REMEDIATION
BIDDEFORD, MAINE**

DWG:

A3

BY:

BMD

DATE:

1/2014

SHEET TITLE:

**MAINE ENERGY
1976 SANBORN MAP**

JN:

12-3240.1

REV:

JKC

SCALE:

NTS

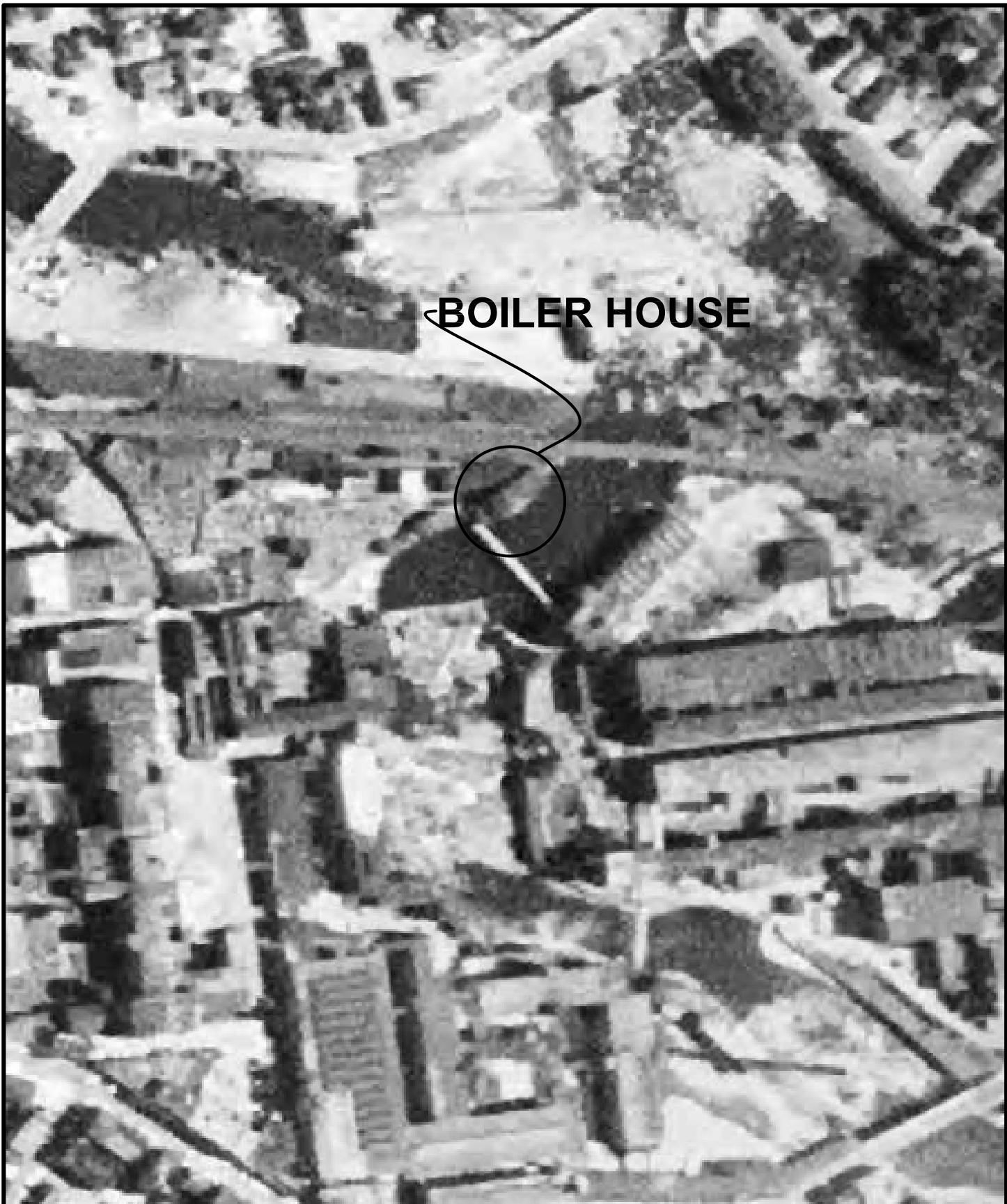
REV DATE: **1/2014**





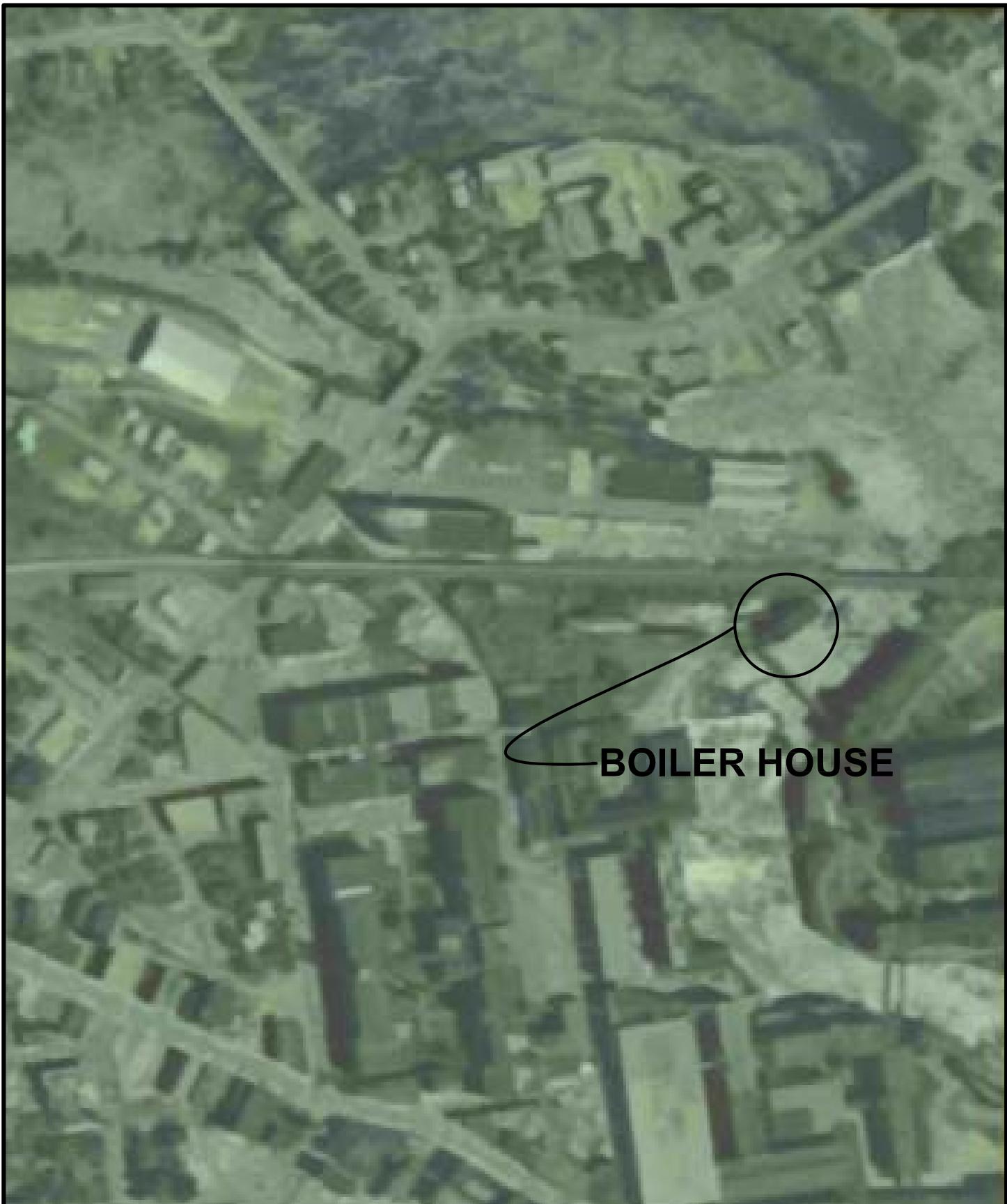
PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE		DWG:	A4	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1956 AERIAL PHOTOGRAPHY		JN:	12-3240.1	DATE:	1/2014
	SCALE:	NTS	REV:	JKC	REV DATE:	1/2014





PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A5	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1960 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





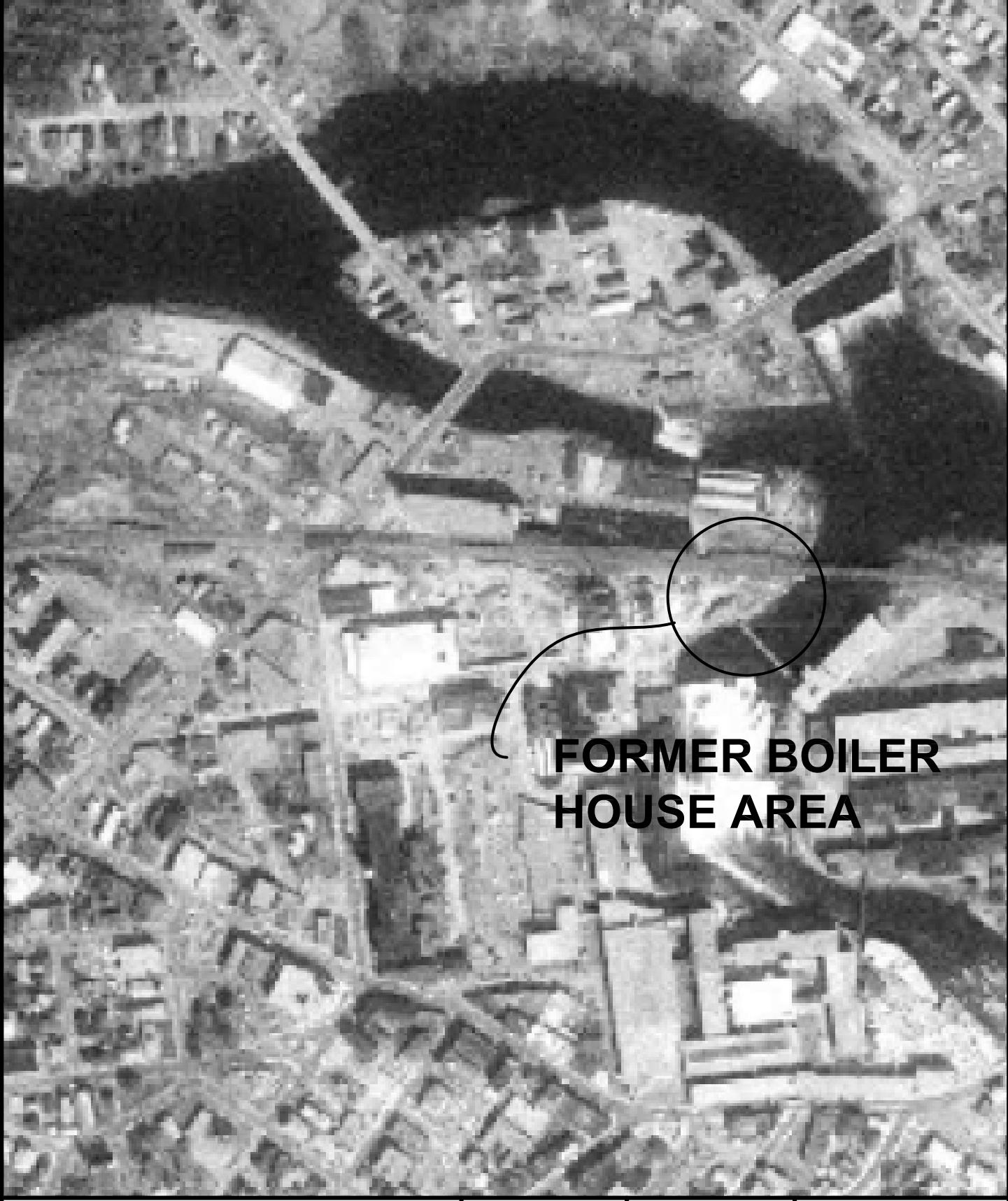
PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A6	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1970 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A7	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1975 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





FORMER BOILER HOUSE AREA

PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A8	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1986 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





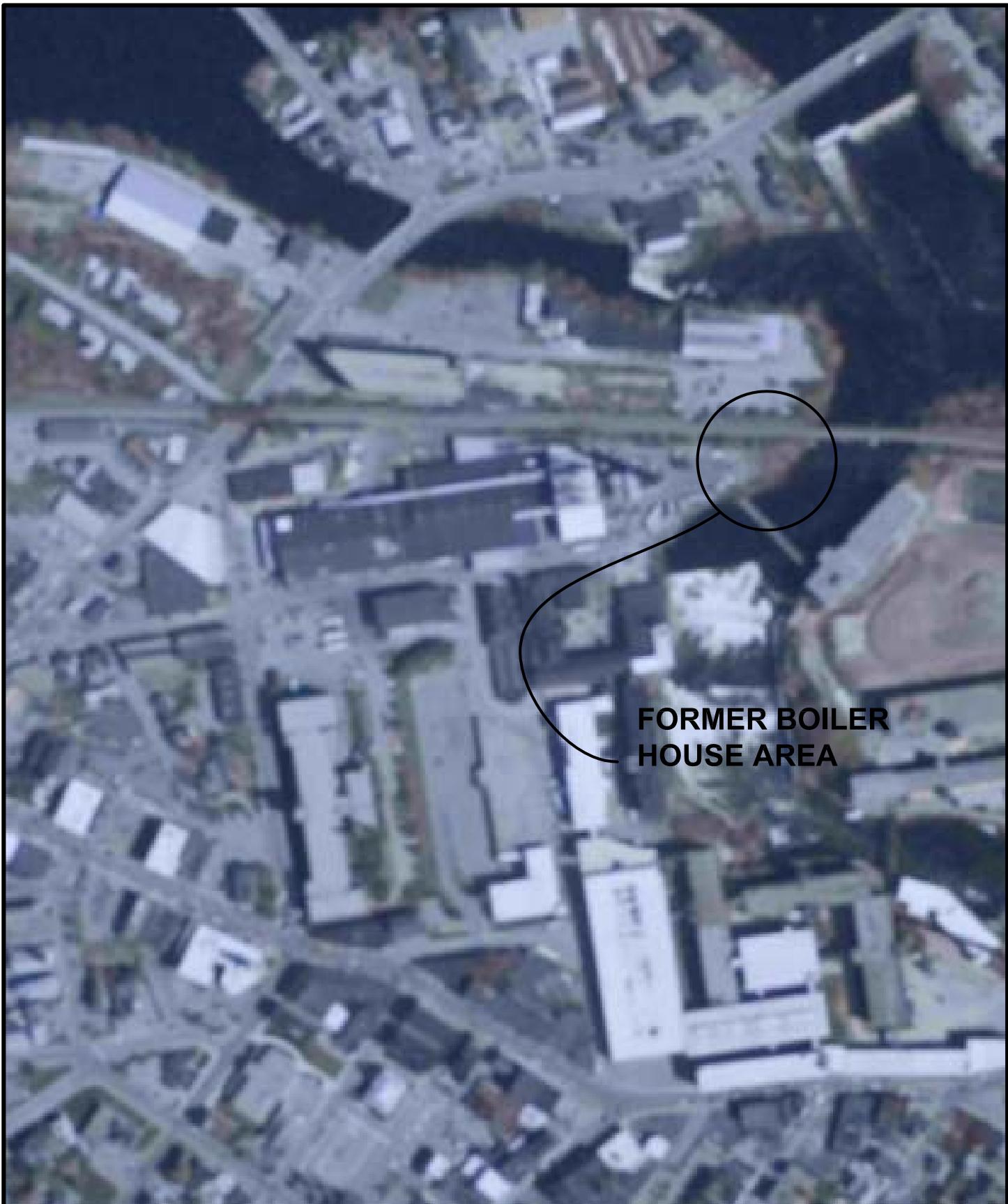
**FORMER BOILER
HOUSE AREA**

PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A9	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1990 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014





PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE		DWG:	A10	BY:	BMD
SHEET TITLE:	MAINE ENERGY 1998 AERIAL PHOTOGRAPHY		JN:	12-3240.1	DATE:	1/2014
	SCALE:	NTS	REV:	JKC	REV DATE:	1/2014
						 The logo for Summit Environmental Consultants features a stylized blue mountain peak graphic above the company name "Summit" in a bold, serif font. Below "Summit" is the word "Environmental Consultants" in a smaller, sans-serif font.



PROJECT TITLE:	PCB REMEDIATION BIDDEFORD, MAINE	DWG:	A11	BY:	BMD
SHEET TITLE:	MAINE ENERGY 2003 AERIAL PHOTOGRAPHY	JN:	12-3240.1	DATE:	1/2014
		SCALE:	NTS	REV:	JKC
				REV DATE:	1/2014



Appendix B

Boring Logs

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-201
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/20	0-2.5		0-3" LOAM 3-20" SAND & GRAVEL Fill (Sample 3-30")		LOAM	
2							SAND & GRAVEL	
	SS-2	30/9	2.5-5'		0-6" Same As Above 6-9" Fill with BRICK (Sample 4-5")		Fill with BRICK	
4								
6								
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Auger refusal @2.5' move to SE-SB-202			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: 3"-4', and 4'-5'			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-202
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/20	0-2.5		0-4" Black SAND & GRAVEL 4-6" ROCK 6-20" Brown Fine to Medium SAND, Trace GRAVEL (Sample at 4"-2.5')		SAND & GRAVEL	
2							ROCK	
4								
6	SS-2	30/12	2.5-5		0-3" Same as above 3-6" ROCK 6-12" Black ASH with GRAVEL (Sample at 4-5') 0-6" Same as above		ROCK	
	SS-3-3	6/6	5-5.5'				ASH w/ GRAVEL	
8					Refusal @ 5.5'			
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 5.5'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: 4"-2.5', and 4'-5'			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-203
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Location: BIDDEFORD, ME				Sheet:			1 of 1	
				Chkd by:			JKC	
Drilling Co:	SUMMIT Geoengineering Services			Boring Location: see location map				
Personnel:	Craig Coolidge			Top of PVC Casing Elevation:				
Summit Staff:	John Cressey			Date started: 1/10/2013 Date Completed: 1/10/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)					SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
	SS-1	30/14	0-2.5'		0-2" Black SAND		SAND	
					2-14" Brown SAND Trace GRAVEL (Fill) (Sample 2"-2.5")		SAND Trace GRAVEL	
2								
	SS-2	30/24	2.5-5'		0-3" ROCK		ROCK	
					3-6" SAND and GRAVEL		SAND & GRAVEL	
4					6-15" Black ASH and GRAVEL (Sample 3'-4')		ASH & GRAVEL	
					15-24" CONCRETE and GRAVEL (Sample 4'-5')		CONCRETE & GRAVEL	
	SS-3	30/26	5-7.5'		0-26" SILTY-SAND, Trace Brick (Sample 5'-7.5")		SILTY SAND	
6								
	SS-4	30/16	7.5-10'		0-6" Same As Above			
8					6-16" ROCK Fragments		ROCK	
10					Refusal @ 10'			
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft	<5%	1. Samples taken at the following intervals: 2"-2.5', 3'-4', 4'-5', 5'-7.5'			
4-10	Loose	2-4	Soft	trace	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30	Compact	4-8	Firm	5-15				
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-204
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	30/24	0-2.5'		0-3" LOAM 3-24" Brown SAND and GRAVEL (Fill) (Sample 3"-3')		LOAM	
4	SS-2	30/22	2.5-5'		0-6" Same As Above 6-22" Black ASH with ROCK (Sample 3'-5')		SAND & GRAVEL	
6	SS-3	30/16	2.5-5'		0-2" Same As Above 2-6" ROCK Fragments Refusal @ 5'		ASH with ROCK	
8								
10								
12								
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition		NOTES:			
Blows/ft. Density	Blows/ft.	Consistency						
0-4 V. Loose	<2	V. soft	<5%	trace	1. Samples taken at the following intervals: 3"-3', 3'-5'			
4-10 Loose	2-4	Soft	5-15	little	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30 Compact	4-8	Firm	15-25	some				
30-50 Dense	8-15	Stiff	>25	and				
>50 V. Dense	15-30	V. Stiff						
	>30	Hard						

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-205
				Project:	MAINE ENERGY PHASE III SUPPORT		Project #:	12-3259.1
				Location:	BIDDEFORD, ME		Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:	SUMMIT Geoengineering Services		Boring Location: see location map					
Personnel:	Craig Coolidge		Top of PVC Casing Elevation:					
Summit Staff:	John Cressey		Date started: 1/10/2013 Date Completed: 1/10/2013					
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)					SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
2	SS-1	30/22	0-2.5'		0-3" LOAM 3-22" Brown Fine SAND and GRAVEL (Fill) (Sample 3"-3.0")		LOAM	
4	SS-2	30/24	2.5-5'		0-4" Same As Above 4-8" ROCK 8-10" ASH 10-24" ROCK and Brown Coarse SAND		SAND & GRAVEL	
6	SS-3	30/5	5-7.5'		0-5" Same As Above (Sample 3.5'-7.5')		ROCK	
8	SS-4	30/14	7.5-10'		0-14" Same As Above		ASH	
10	SS-5	30/26	10-12.5'		0-3" Same As Above 3-26" Brownish-Gray SILTY-CLAY moist with Trace GRAVEL (Sample 10.5-12")		ROCK & SAND	
12							SILTY-CLAY	
14					Cease @ 12.5'			
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Samples taken at the following intervals: 3"-3', 3.5'-7', 10.5-12.5'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30	Compact	4-8	Firm	5-15 little				
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-206
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/26	0-2.5'		0-3" LOAM 3-16" Brown SAND and GRAVEL 16-26" ROCK Fragments, Trace SAND (Sample 3"-2")		LOAM SAND & GRAVEL	
2					0-10" Same As Above			
4					0-8" Same As Above			
6	SS-3	30/8	5-7.5'		0-8" Same As Above		ROCK, Trace SAND	
8	SS-4	30/8	7.5-10'		0-8" Same As Above			
10	SS-5	30/18	10-12.5'		0-12" Same As Above 12-18" ROCK (Sample 2'-12')		ROCK	
12					Cease @ 12.5'			
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Samples taken at the following intervals: 3"-2', 2'-12'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30	Compact	4-8	Firm	5-15 little				
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-207
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Location: BIDDEFORD, ME							Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:	SUMMIT Geoengineering Services			Boring Location: see location map				
Personnel:	Craig Coolidge			Top of PVC Casing Elevation:				
Summit Staff:	John Cressey			Date started: 1/10/2013 Date Completed: 1/10/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)					SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
	SS-1	30/17	0-2.5'		0-3" LOAM 3-17" Black SAND & GRAVEL (Fill)		LOAM	
2					(Sample 3"-3')		SAND & GRAVEL	
	SS-2	30/20	2.5-5'				ASH & ROCK	
4							ROCK	
6	SS-3	24/16	5-7'		0-6" Same As Above 6-18" Black ASH and ROCK 18-20" ROCK		SAND & GRAVEL	
					(Sample 3'-4.5')		ROCK	
8					0-10" Brown SAND and GRAVEL 10-16" ROCK			
10								
12					(Sample 5'-6')			
14								
16								
18								
20					Refusal @ 7'			
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Samples taken at the following intervals: 3"-3', 3'-4.5', 5'-6'			
4-10	Loose	2-4	Soft	<5%	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30	Compact	4-8	Firm	trace				
30-50	Dense	8-15	Stiff	5-15				
>50	V. Dense	15-30	V. Stiff	15-25				
		>30	Hard	>25				
				and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-208
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/20	0-2.5'		0-3" LOAM 3-20" Brown SAND & GRAVEL (Sample 3"-4")		LOAM	
2							SAND & GRAVEL	
	SS-2	30/4	2.5'-5'		0-1" Same As Above 1-4" BRICK and SAND (Fill) (Sample 4'-6")		BRICK & SAND	
4							SAND & GRAVEL	
	SS-3	24/16			0-12" Sam As Above 12-16" Gray SAND and GRAVEL (Sample 6'-8")		SILTY-SAND	
6								
	SS-4	30/20	7.5-10'		0-6" Same As Above 6-20" Brown SILTY-SAND, Moist (Sample 8'-10")			
8								
					Refusal @ 10"			
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Samples taken at the following intervals: 3"-4", 4'-6", 6'-8", 8-10'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
10-30	Compact	4-8	Firm	5-15 little				
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-209
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/20	0-2.5'		0-2" CONCRETE 2-20" Brown SAND & GRAVEL (Fill)		CONCRETE	
2							SAND & GRAVEL	
	SS-2	30/18	2.5'-5'		0-3" Same As Above 3-18" Black SAND and BRICK			
4							SAND & BRICK	
	SS-3	30/16`5-7.5'			0-16" Same As Above			
6								
	SS-4	30/20	7.5-10'		0-15" Same As Above 15-20" ASH (Clinker)			
8								
	SS-5	30/16	10-12.5		0-6" Same as Above		ASH	
10								
					(Sample 3'-9')			
12					Cease @ 12.5'			
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Voids noted from 9-10'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: 2"-3', 3'-9', 9'-12.5'			
10-30	Compact	4-8	Firm	5-15	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #: SE-SB-210
				Project: <u>MAINE ENERGY PHASE III SUPPORT</u>		Project #: <u>12-3259.1</u>	
				Location: <u>BIDDEFORD, ME</u>		Sheet: <u>1 of 1</u> Chkd by: <u>JKC</u>	
Drilling Co: <u>SUMMIT Geoengineering Services</u>				Boring Location: see location map			
Personnel: <u>Craig Coolidge</u>				Top of PVC Casing Elevation:			
Summit Staff: <u>John Cressey</u>				Date started: <u>1/10/2013</u> Date Completed: <u>1/10/2013</u>			
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH			
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)
Model:	9500-VTR	Hammer:	N/A	See Report			
Method:	Direct Push	Fall:	N/A				
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)	
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.			
SS-1	20/10	0-1.75'		0-20" Black SILTY-SAND (Sample 0-20")	SILTY-SAND		
2				Refusal @ 20"			
4							
6							
8							
10							
12							
14							
16							
18							
20							
Granular Soils	Cohesive Soils		% Composition	NOTES:			
Blows/ft. Density	Blows/ft.	Consistency		1. Boring reached refusal at 20" in three different locations 2. Samples taken at the following intervals: 0-20" 2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
0-4 V. Loose	<2	V. soft	<5% trace				
4-10 Loose	2-4	Soft	5-15 little				
10-30 Compact	4-8	Firm	15-25 some				
30-50 Dense	8-15	Stiff	>25 and				
>50 V. Dense	15-30	V. Stiff					
	>30	Hard					

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-211
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE III SUPPORT			Project #:	12-3259.1
				Location: BIDDEFORD, ME			Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map				
Personnel: Craig Coolidge				Top of PVC Casing Elevation:				
Summit Staff: John Cressey				Date started: 1/10/2013 Date Completed: 1/10/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	N/A	See Report				
Method:	Direct Push	Fall:	N/A					
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)		
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
SS-1	30/22	0-2.5'		0-2" LOAM 2-22" Brown SAND & GRAVEL (Fill)	(Sample 2"-3.5")	LOAM		
SS-2	30/20	2.5-5'		0-16" Same As Above		SAND & GRAVEL		
SS-3	30/10	5-7.5'		16-20" Dark Brown SAND and BRICK				
SS-4	18/5	7.5-9'		0-10" Same As Above 0-5" Same As Above	(Sample 4-9")	SAND & BRICK		
				Refusal @ 9'				
10								
12								
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition	NOTES:				
Blows/ft.	Density	Blows/ft.	Consistency	1. Samples taken at the following intervals: 2"-3.5', 4'-9' 2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
0-4	V. Loose	<2	V. soft	<5%	trace			
4-10	Loose	2-4	Soft	5-15	little			
10-30	Compact	4-8	Firm	15-25	some			
30-50	Dense	8-15	Stiff	>25	and			
>50	V. Dense	15-30	V. Stiff					
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #: SE-SB-212
				Project: <u>MAINE ENERGY PHASE III SUPPORT</u>		Project #: <u>12-3259.1</u>	
				Location: <u>BIDDEFORD, ME</u>		Sheet: <u>1 of 1</u> Chkd by: <u>JKC</u>	
Drilling Co: <u>SUMMIT Geoengineering Services</u>				Boring Location: see location map			
Personnel: <u>Craig Coolidge</u>				Top of PVC Casing Elevation:			
Summit Staff: <u>John Cressey</u>				Date started: <u>1/10/2013</u> Date Completed: <u>1/10/2013</u>			
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH			
Vehicle:	AMS	Type:	Split Spoon	Date	Depth (ft)	Reference	Groundwater Elevation (ft)
Model:	9500-VTR	Hammer:	N/A	See Report			
Method:	Direct Push	Fall:	N/A				
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)	
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.			
SS-1	20/15	0-20"		0-2" LOAM 2-15" Brown SAND and GRAVEL	LOAM		
2				Refusal @ 20"			
4							
6							
8							
10							
12							
14							
16							
18							
20							
Granular Soils	Cohesive Soils		% Composition		NOTES:		
Blows/ft.	Density	Blows/ft.	Consistency				
0-4	V. Loose	<2	V. soft	<5%	trace	1. Boring reached refusal at 20" in three locations	
4-10	Loose	2-4	Soft	5-15	little	2. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)	
10-30	Compact	4-8	Firm	15-25	some		
30-50	Dense	8-15	Stiff	>25	and		
>50	V. Dense	15-30	V. Stiff				
		>30	Hard				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-213
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1
				Location:	BIDDEFORD, ME		Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:	SUMMIT Geoengineering Services		Boring Location: see location map					
Personnel:	Craig Coolidge		Top of PVC Casing Elevation:					
Summit Staff:	John Cressey		Date started: 2/19/2013 Date Completed: 2/19/2013					
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/15	0-2.5'		0-3" Asphalt 3-15" Brown SAND & GRAVEL (Fill)		ASPHALT	
2							SAND & GRAVEL	
	SS-2	30/30	2.5-5'		0-6" Blackish-Brown MEDIUM SAND 6-15" Brown MEDIUM SAND & GRAVEL 15-18" ROCK 18-28" Brown SAND & GRAVEL 28-30" BRICK (Demo Debris)		SAND SAND & GRAVEL ROCK SAND & GRAVEL DEMO DEBRIS	
4								
	SS-3	30/30	5-7.5'		0-10" Brown MEDIUM SAND 10-14" ASH & BRICK 14-18" ROCK 18-22" CONCRETE 22-30" BRICK & SAND		SAND DEMO DEBRIS ROCK DEMO DEBRIS	
6								
8					REFUSAL @ 7.5'			
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 7.5'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-7.5')			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-214
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	30/23	0-2.5'		0-4" LOAM & ORGANIC MATERIAL 4-19" Brown SAND & GRAVEL (Fill) 19-23" Reddish-Brown FINE SAND		LOAM SAND & GRAVEL SAND	
							SAND & GRAVEL	
							ROCK	
	SS-2	30/21	2.5-5'		5-8" ROCK & SANDY GRAVEL 8-14" Black ASH & ROCK 14-20" ROCK 20-21" ASH		ROCK SAND GRAVEL DEMO DEBRIS ROCK DEMO DEBRIS	
4								
	SS-3	24/24	5-7.1'		0-12" Black ASH, SAND & GRAVEL 12-14" BRICK w/ Sand & Ash 14-16" CONCRETE 16-24" BRICK w/ Sand & Ash		DEMO DEBRIS	
6								
					REFUSAL @ 7.1'			
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition		NOTES:		
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft	<5%	trace	1. Refusal @ 7.1'		
4-10	Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-7')		
10-30	Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)		
30-50	Dense	8-15	Stiff	>25	and			
>50	V. Dense	15-30	V. Stiff					
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-215		
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1		
				Location:	BIDDEFORD, ME		Sheet:	1 of 1		
							Chkd by:	JKC		
Drilling Co:	SUMMIT Geoengineering Services			Boring Location: see location map						
Personnel:	Craig Coolidge			Top of PVC Casing Elevation:						
Summit Staff:	John Cressey			Date started: 2/19/2013 Date Completed: 2/19/2013						
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH						
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model:	9500-VTR	Hammer:	140LR	See Report						
Method:	H.S.A.	Fall:	30"							
Depth (ft.)					SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)		
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.						
2	SS-1	30/20	0-2.5'		0-3" ASPHALT		ASPHALT			
					3-15" Brown SAND & GRAVEL (Fill)		SAND & GRAVEL			
					15-16" ROCK		ROCK			
					16-20" Brown SAND & GRAVEL (Fill)		SAND & GRAVEL			
4	SS-2	30/26	2.5-5'		0-2" Brown SAND & GRAVEL (Fill)		DEMO DEBRIS			
					2-8" Black ASH & SILTY SAND		SAND & GRAVEL			
					8-15" Brown SAND & GRAVEL		DEMO DEBRIS			
					15-17" BRICK		SAND & GRAVEL			
6	SS-3	30/28	5-7.5'		17-26" Brown SAND & GRAVEL (Fill) w Ash & Rock		ROCK			
					0-3" Brown SAND & GRAVEL (Fill) w/ Ash & Rock		DEMO DEBRIS			
					3-4" ROCK		SAND & GRAVEL			
					4-7" BRICK with Ash & Sand		ROCK			
8	SS-4	30/15	7.5-10'		7-9" CONCRETE		DEMO DEBRIS			
					9-16" Brownish Red SAND & GRAVEL		SAND & GRAVEL			
					16-28" BRICK with Ash & Sand		DEMO DEBRIS			
					0-3" BRICK with Ash & Sand					
					3-15" Black ASH trace brick					
10					REFUSAL @ 10'					
12										
14										
16										
18										
20										
Granular Soils		Cohesive Soils		% Composition						
Blows/ft.	Density	Blows/ft.	Consistency	NOTES:						
0-4	V. Loose	<2	V. soft	<5%	trace	1. Refusal @ 10'				
4-10	Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')				
10-30	Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
30-50	Dense	8-15	Stiff	>25	and					
>50	V. Dense	15-30	V. Stiff	>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-216
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/16	0-2.5'		0-2.5" ASPHALT 2.5-24" Brown SAND & GRAVEL (Fill)		ASPHALT	
2							SAND & GRAVEL	
	SS-2	30/28	2.5-5'		0-3" Brown SAND & GRAVEL (Fill) 3-10" Black SAND (Fill)		SAND	
4					10-14" BRICK			
	SS-3	30/16	5-6.8'		14-20" CONCRETE (Fill) 20-28" CONCRETE & SAND (Fill)		DEMO DEBRIS	
6					REFUSAL @ 6.8'			
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 10'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (2.5-24") (2-4') (4-6')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-217
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD				ESTIMATED GROUND WATER DEPTH				
Vehicle: AMS	Type: 24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model: 9500-VTR	Hammer: 140LR	See Report						
Method: H.S.A.	Fall: 30"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	30/12	0-2.5'		0-10" Black SAND, trace gravel 10-12" Black ORGANIC MATERIAL		SAND	
4	SS-2	30/16	2.5-5'		0-10" Black ORGANIC MATERIAL 10-12" Brown CLAY 12-16" Brown SAND & GRAVEL		LOAM	
6	SS-3	30/0	5-7.5'		NO RECOVERY		CLAY	
8	SS-4	30/3	7.5-10		0-3" Black ASH & SAND		SAND & GRAVEL	
10					CEASE @ 10' (HOLE COLLAPSED)		DEMO DEBRIS	
12								
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition		NOTES:			
Blows/ft. Density	Blows/ft.	Consistency						
0-4 V. Loose	<2	V. soft	<5%	trace	1. Cease @ 10'			
4-10 Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(8-10')			
10-30 Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50 Dense	8-15	Stiff	>25	and				
>50 V. Dense	15-30	V. Stiff						
	>30	Hard						

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-218
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1
				Location:	BIDDEFORD, ME		Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:	SUMMIT Geoengineering Services		Boring Location: see location map					
Personnel:	Craig Coolidge		Top of PVC Casing Elevation:					
Summit Staff:	John Cressey		Date started: 2/19/2013 Date Completed: 2/19/2013					
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	24/18	0-2'		0-3" LOAM		LOAM	
					3-10" Brown SAND & GRAVEL		SAND & GRAVEL	
2					10-18" Brown FINE SAND		SAND	
	SS-2	24/20	2-4"		0-10" Brown FINE SAND			
					10-20" Gray SILTY CLAY		CLAY	
4					CEASE @ 4'			
6								
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Cease @ 4'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-219
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	30/16	0-2.5'		0-3" LOAM & ORGANIC MATERIAL 3-6" Dark brown SAND & GRAVEL 6-14" Brown FINE TO MEDIUM SAND, trace gravel 14-16" CONCRETE		LOAM	
4	SS-2	30/16	2.5-5'		0-10" SAND AND GRAVEL 10-14" WOOD 14-16" SAND & GRAVEL		SAND & GRAVEL	
6	SS-3	30/5	5-7.5'		0-5" GRAVEL		DEMO DEBRIS	
8	SS-4	30/10	7.5-10'		0-10" GRAVEL		SAND & GRAVEL	
10					REFUSAL @ 10'		GRAVEL	
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 10'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-220
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/15	0-2.5'		0-4" LOAM 4-15" Brown SAND & GRAVEL (Fill)		LOAM	
2							SAND & GRAVEL	
	SS-2	30/16	2.5-5'		0-2" Brown SAND & GRAVEL (Fill) 2-4" ASH 4-8" Brown SAND & BRICK 8-10" ROCK, 10-15" ASH & BRICK 15-16" CONCRETE		DEMO DEBRIS	
4							SAND, DEBRIS	
	SS-3	30/16	5-7.5'		0-6" ASH & BRICK 6-16" Black ASH, trace sand		ROCK, DEBRIS	
6							DEMO DEBRIS	
	SS-4	30/16	7.5-10'		0-16" Black ASH, trace sand		DEBRIS, SAND	
8								
10	SS-5	12/8	10-11'		0-3" Black ASH, trace sand 3-8" ROCK		ROCK	
12					REFUSAL @ 11'			
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 11'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')(10-11')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-221
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/16	0-2.5'		0-8" Brown SAND & GRAVEL 8-16" BRICK		SAND & GRAVEL	
2							DEMO DEBRIS	
	SS-2	30/30	2.5-5'		0-4" BRICK 4-30" BRICK, CONCRETE, Some Sand		DEBRIS, SAND	
4								
6					REFUSAL @ 5'			
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 5'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: (0-2')(2-4')(4-5')			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-222
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Location: BIDDEFORD, ME				Sheet: 1 of 1			Chkd by:	JKC
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map				
Personnel: Craig Coolidge				Top of PVC Casing Elevation:				
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	concrete punch	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	24"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	SAMPLE DESCRIPTION			Stratum	Field Screening (ppmv)
				Blows/6 in.				
	SS-1	24/20	0-2'	0-3" LOAM			LOAM	
				3-15" SAND & GRAVEL			SAND & GRAVEL	
				15-16" Black ASH			DEMO DEBRIS	
2	SS-2	24/20	2-4'	16-20" Brown SAND & GRAVEL			SAND & GRAVEL	
				0-10" Brown SAND & GRAVEL			DEMO DEBRIS	
				10-14" BRICK			SAND	
4	SS-3	24/18	4-6'	14-20" Brown MEDIUM SAND				
				0-18" Brown MEDIUM SAND w/ Rock & Brick				
6	SS-4	24/16	6-8'	0-16" Brown MEDIUM SAND w/ Rock, Brick, Clinker			SAND, DEBRIS	
				2-18" Brown FINE SAND			SAND	
8	SS-5	24/18	8-10'	CEASE @ 10'				
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition				
Blows/ft.	Density	Blows/ft.	Consistency	NOTES:				
0-4	V. Loose	<2	V. soft	<5%	trace		1. Cease @ 10'	
4-10	Loose	2-4	Soft	5-15	little		2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')	
10-30	Compact	4-8	Firm	15-25	some		3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)	
30-50	Dense	8-15	Stiff	>25	and			
>50	V. Dense	15-30	V. Stiff					
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-223
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	concrete punch	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	24"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	24/18	0-2'		0-3" LOAM 3-18" Brown SAND & GRAVEL		LOAM	
2							SAND & GRAVEL	
	SS-2	24/18	2-4'		0-10" Brown SAND & GRAVEL 10-12" BRICK 12-18" BRICK & SAND		DEMO DEBRIS	
4							DEBRIS & SAND	
	SS-3	24/20	4-6'		0-15" BRICK & SAND 15-20" Brown SAND w/ Brick			
6								
	SS-4	24/16	6-8'		0-8" Brown SAND w/ Brick 8-16" Brown MEDIUM SAND			
8								
	SS-5	24/2	8-10'		0-2" Brown MEDIUM SAND		SAND	
10					CEASE @ 10'			
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Cease @ 10'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-224		
Drilling Co:	SUMMIT Geoengineering Services			Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1		
Personnel:	Craig Coolidge			Location:	BIDDEFORD, ME		Sheet:	1 of 1		
Summit Staff:	John Cressey						Chkd by:	JKC		
DRILLING METHOD		SAMPLER		Boring Location: see location map						
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	ESTIMATED GROUND WATER DEPTH			
Model:	9500-VTR	Hammer:	140LR	See Report			Groundwater Elevation (ft)			
Method:	H.S.A.	Fall:	30"							
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)				
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.						
2	SS-1	30/22	0-2.5'	0-3" LOAM & ORGANIC MATERIAL			LOAM			
				3-5" SAND & GRAVEL			SAND & GRAVEL			
4				5-12" Black ASH & SAND			DEMO DEBRIS			
				12-14" Brown MEDIUM SAND, 14-18" Black ASH			DEBRIS & SAND			
6	SS-2	30/10	2.5-5'	18-22" Reddish-brown SAND			SAND			
				0-10" Reddish-brown SAND w/ Gravel						
8	SS-3	30/0	5-7.5'	NO RECOVERY						
	SS-4	12/6	7.5-8'	0-6" GRAVEL			GRAVEL			
10				REFUSAL @ 8'						
12										
14										
16										
18										
20										
Granular Soils		Cohesive Soils		% Composition						
Blows/ft.	Density	Blows/ft.	Consistency	NOTES:						
0-4	V. Loose	<2	V. soft	<5%	trace	1. Refusal @ 8'				
4-10	Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')(6-8')				
10-30	Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
30-50	Dense	8-15	Stiff	>25	and					
>50	V. Dense	15-30	V. Stiff							
		>30	Hard							

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-225
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	30/8	0-2.5'		0-3" LOAM & ORGANIC MATERIAL 3-5" Brown FINE SAND 5-8" Brown SAND & GRAVEL		LOAM	
2							SAND	
	SS-2	30/10	2.5-5'		0-10" Brown SAND & GRAVEL			
4								
	SS-3	30/12	5-7.5'		0-12" Brown SAND & GRAVEL			
6								
8					REFUSAL @ 7.5'			
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 7.5"			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-7.5')			
10-30	Compact	4-8	Firm	trace				
30-50	Dense	8-15	Stiff	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
>50	V. Dense	15-30	V. Stiff	15-25				
		>30	Hard	>25				
				and				

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-226
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
				Location: BIDDEFORD, ME			Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map				
Personnel: Craig Coolidge				Top of PVC Casing Elevation:				
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)		
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
SS-1	30/10	0-2.5'		0-3" LOAM 3-10" Brown SAND & GRAVEL	LOAM			
2					SAND & GRAVEL			
SS-2	30/2	2.5-5'		0-2" Brown SILTY SAND	SAND			
4								
6				REFUSAL @ 5'				
8								
10								
12								
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition	NOTES:				
Blows/ft.	Density	Blows/ft.	Consistency	1. Refusal @ 5' 2. Samples taken at the following intervals: (0-2')(2-4')(4-5') 3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
0-4	V. Loose	<2	V. soft	<5%	trace			
4-10	Loose	2-4	Soft	5-15	little			
10-30	Compact	4-8	Firm	15-25	some			
30-50	Dense	8-15	Stiff	>25	and			
>50	V. Dense	15-30	V. Stiff					
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-227
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1
				Location:	BIDDEFORD, ME		Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:	SUMMIT Geoengineering Services		Boring Location: see location map					
Personnel:	Craig Coolidge		Top of PVC Casing Elevation:					
Summit Staff:	John Cressey		Date started: 2/20/2013 Date Completed: 2/20/2013					
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	24"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	24/18	0-2'		0-3" LOAM		LOAM	
					3-10" SAND & GRAVEL		SAND & GRAVEL	
2					10-18" Brown MEDIUM SAND		SAND	
	SS-2	24/20	2-4'		0-20" BRICK			
					0-20" BRICK & CONCRETE		DEMO DEBRIS	
4	SS-3	24/20	4-6'		0-10" BRICK & CONCRETE			
					10-14" Brown SILTY SAND		SAND	
6	SS-4	24/18	6-8'		14-18" ROCK		ROCK	
					REFUSAL @ 8'			
8								
10								
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Refusal @ 8'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-228		
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1		
				Location:	BIDDEFORD, ME		Sheet:	1 of 1		
							Chkd by:	JKC		
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map						
Personnel: Craig Coolidge				Top of PVC Casing Elevation:						
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013						
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH						
Vehicle:	AMS	Type:	Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model:	9500-VTR	Hammer:	140LR	See Report						
Method:	H.S.A.	Fall:	24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	SAMPLE DESCRIPTION			Stratum	Field Screening (ppmv)		
				Blows/6 in.						
	SS-1	24/18	0-2'	0-3" LOAM			LOAM			
				3-10" Brown SAND & GRAVEL			SAND & GRAVEL			
2				10-18" Brown SAND AND GRAVEL w/ Rock			SAND & GRAVEL			
	SS-2	24/20	2-4'	0-20" BRICK & CONCRETE						
				0-10" BRICK & CONCRETE			DEMO DEBRIS			
4	SS-3	24/20	4-6'	10-20" Black ASH & BRICK						
				0-18" BRICK & SAND			DEBRIS, SAND			
6	SS-4	24/18	6-8'	0-10" BRICK & SAND						
				10-14" Brown SILTY SAND			SAND			
8	SS-5	24/16	8-10'	14-16" ROCK			ROCK			
10				CEASE @ 10'						
12										
14										
16										
18										
20										
Granular Soils		Cohesive Soils		% Composition						
Blows/ft.	Density	Blows/ft.	Consistency	NOTES:						
0-4	V. Loose	<2	V. soft	<5%	trace					
4-10	Loose	2-4	Soft	5-15	little					
10-30	Compact	4-8	Firm	15-25	some					
30-50	Dense	8-15	Stiff	>25	and					
>50	V. Dense	15-30	V. Stiff							
		>30	Hard							

1. Cease @ 10'
 2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')
 3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-229
				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1
				Location:	BIDDEFORD, ME		Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co:		SUMMIT Geoengineering Services		Boring Location: see location map				
Personnel:		Craig Coolidge		Top of PVC Casing Elevation:				
Summit Staff:		John Cressey		Date started: 2/20/2013 Date Completed: 2/20/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	24"					
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	24/20	0-2'		0-3" LOAM		LOAM	
					3-8" Brown SAND & GRAVEL		SAND & GRAVEL	
					8-20" BRICK & CONCRETE, trace sand		DEMO DEBRIS	
2	SS-2	24/20	2-4'		0-20" Brown SAND & GRAVEL, trace brick			
4	SS-3	24/20	4-6'		0-10" Brown SAND & GRAVEL, trace brick			
					10-20" BRICK, ASH, CONCRETE			
6	SS-4	24/18	6-8'		0-18" Brown SAND & GRAVEL wet			
8	SS-5	24/20	8-10'		0-20" Brown SAND & GRAVEL wet			
10					CEASE @ 10'			
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Cease @ 10'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #: SE-SB-230
				Project: <u>MAINE ENERGY PHASE II SUPPORT</u>		Project #: 12-3259.1	
				Location: <u>BIDDEFORD, ME</u>		Sheet: 1 of 1	
						Chkd by: JKC	
Drilling Co: <u>SUMMIT Geoengineering Services</u>				Boring Location: see location map			
Personnel: <u>Craig Coolidge</u>				Top of PVC Casing Elevation:			
Summit Staff: <u>John Cressey</u>				Date started: <u>2/20/2013</u> Date Completed: <u>2/20/2013</u>			
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH			
Vehicle:	AMS	Type:	Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)
Model:	9500-VTR	Hammer:	140LR	See Report			
Method:	H.S.A.	Fall:	24"				
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)	
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.			
2	SS-1	24/20	0-2'		0-3" LOAM 3-20" Brown SAND & GRAVEL With Rock	LOAM	
4	SS-2	20/10	2-4'		0-10" Brown SAND & GRAVEL With Rock	SAND & GRAVEL	
6					REFUSAL @ 4'		
8							
10							
12							
14							
16							
18							
20							
Granular Soils	Cohesive Soils		% Composition		NOTES:		
Blows/ft.	Density	Blows/ft.	Consistency				
0-4	V. Loose	<2	V. soft	<5%	trace	1. Refusal @ 4'	
4-10	Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')	
10-30	Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)	
30-50	Dense	8-15	Stiff	>25	and		
>50	V. Dense	15-30	V. Stiff				
		>30	Hard				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-231
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle: AMS	Type: Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model: 9500-VTR	Hammer: 140LR	See Report						
Method: H.S.A.	Fall: 24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	24/20	0-2'		0-3" LOAM 3-8" SAND & GRAVEL (Fill) 8-20" BRICK, ASPHALT, CONCRETE		LOAM SAND & GRAVEL	
					0-15" BRICK, ASPHALT, CONCRETE 15-22" ROCK		DEMO DEBRIS	
4	SS-2	24/22	2-4'				ROCK	
					0-10" BRICK & CLINKER 10-24" CONCRETE w/ trace sand & gravel		DEMO DEBRIS	
6	SS-3	24/24	4-6'				DEBRIS, SAND	
					0-6" CONCRETE w/ trace sand & gravel 6-20" Reddish-brown SAND & GRAVEL			
8	SS-4	24/20	6-8'					
					0-20" Reddish-brown SAND & GRAVEL		SAND & GRAVEL	
10	SS-5	24/20	8-10'					
					CEASE @ 10'			
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Cease @ 10'			
4-10	Loose	2-4	Soft	<5% trace	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')			
10-30	Compact	4-8	Firm	5-15 little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25 some				
>50	V. Dense	15-30	V. Stiff	>25 and				
		>30	Hard					

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-232
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE II SUPPORT		Project #: 12-3259.1		
				Location: BIDDEFORD, ME		Sheet: 1 of 1 Chkd by: JKC		
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map				
Personnel: Craig Coolidge				Top of PVC Casing Elevation:				
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	24"					
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)		
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
SS-1	24/6	0-2'		0-3" LOAM 3-5" SAND & GRAVEL (Fill) 5-6" BRICK w/ Concrete	LOAM	SAND & GRAVEL DEMO DEBRIS		
2				REFUSAL @ 2'				
4								
6								
8								
10								
12								
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition	NOTES:				
Blows/ft.	Density	Blows/ft.	Consistency	1. Refusal @ 2' 2. Samples taken at the following intervals: (0-2') 3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
0-4	V. Loose	<2	V. soft	<5%	trace			
4-10	Loose	2-4	Soft	5-15	little			
10-30	Compact	4-8	Firm	15-25	some			
30-50	Dense	8-15	Stiff	>25	and			
>50	V. Dense	15-30	V. Stiff					
		>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-232A
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle: AMS	Type: Concrete Punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model: 9500-VTR	Hammer: 140LR	See Report						
Method: H.S.A.	Fall: 24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
	SS-1	24/0	0-2'		NO SAMPLE		LOAM SAND & GRAVEL	
2					0-20" CONCRETE & CLINKER w/ Ash			
4	SS-2	24/20	2-4'		0-20" CONCRETE & CLINKER w/ Ash & Brick		DEMO DEBRIS	
6					0-10" CONCRETE & CLINKER w/ Ash & Brick 10-20" Brown SILTY CLAY, trace clay			
8	SS-3	24/20	4-6'		0-20" Reddish-brown MEDIUM SAND & GRAVEL, trace clinker		CLAY	
10	SS-4	24/20	6-8'				DEBRIS, SAND, GRAVEL	
12					CEASE @ 10'			
14								
16								
18								
20								
Granular Soils	Cohesive Soils		% Composition		NOTES:			
Blows/ft. Density	Blows/ft.	Consistency						
0-4 V. Loose	<2	V. soft			1. Cease @ 10'			
4-10 Loose	2-4	Soft	<5%	trace	2. Samples taken at the following intervals: (2-4')(4-6')(6-8')(8-10')			
10-30 Compact	4-8	Firm	5-15	little	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50 Dense	8-15	Stiff	15-25	some				
>50 V. Dense	15-30	V. Stiff	>25	and				
	>30	Hard						

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-233
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
				Location: BIDDEFORD, ME			Sheet:	1 of 1
							Chkd by:	JKC
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map				
Personnel: Craig Coolidge				Top of PVC Casing Elevation:				
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013				
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle:	AMS	Type:	24" SPT	Date	Depth (ft)	Reference	Groundwater Elevation (ft)	
Model:	9500-VTR	Hammer:	140LR	See Report				
Method:	H.S.A.	Fall:	30"					
Depth (ft.)	SAMPLE DESCRIPTION				Stratum	Field Screening (ppmv)		
	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.				
2	SS-1	30/28	0-2.5'	0-3" LOAM 3-6" GRAVEL 6-28" CONCRETE & BRICK, trace clinker	LOAM			
4	SS-2	24/20	2.5-4'	0-2" CLINKER 2-20" Reddish-brown SAND & GRAVEL, trace rock	GRAVEL			
6	SS-3	24/22	4-6'	0-22" Reddish-brown SAND & GRAVEL, trace rock	DEMO DEBRIS			
8	SS-4	24/24	6-8'	0-10" Reddish-brown SAND & GRAVEL, trace rock 10-12" CONCRETE 12-24" Brown SILTY CLAY, trace rock	SAND, GRAVEL, ROCK			
10				REFUSAL @ 8'	DEMO DEBRIS			
12					CLAY, ROCK			
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition			NOTES:	
Blows/ft.	Density	Blows/ft.	Consistency	<2	V. soft	<5%	trace	
0-4	V. Loose	2-4	Soft	5-15	little			1. Refusal @ 8'
4-10	Loose	4-8	Firm	15-25	some			2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')
10-30	Compact	8-15	Stiff	>25	and			3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)
30-50	Dense	15-30	V. Stiff					
>50	V. Dense	>30	Hard					

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-234
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1
Summit Staff: John Cressey							Chkd by:	JKC
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH				
Vehicle: AMS	Type: Concrete punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model: 9500-VTR	Hammer: 140LR	See Report						
Method: H.S.A.	Fall: 24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)
2	SS-1	24/22	0-2'		0-3" LOAM 3-10" SAND & GRAVEL, trace ash & clinker 10-22" Brown SILTY-SAND		LOAM	
					0-2" ROCK 2-10" Black ASH & CLINKER 10-22" Reddish-brown SAND & GRAVEL		SAND	
4	SS-2	24/22	2-4'				ROCK	
							DEMO DEBRIS	
6	SS-3	24/20	4-6'		0-20" Reddish-brown SAND & GRAVEL		SAND & GRAVEL	
					0-10" Reddish-brown SAND & GRAVEL 10-24" Reddish-brown SAND & GRAVEL w/ Rock			
8	SS-4	24/24	6-8'					
					0-20" Reddish-brown SAND & GRAVEL w/ Rock		SAND, GRAVEL, ROCK	
10	SS-5	24/20	8-10'					
					Cease @ 10'			
12								
14								
16								
18								
20								
Granular Soils		Cohesive Soils		% Composition	NOTES:			
Blows/ft.	Density	Blows/ft.	Consistency					
0-4	V. Loose	<2	V. soft		1. Cease @ 10'			
4-10	Loose	2-4	Soft	<5%	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')			
10-30	Compact	4-8	Firm	5-15	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)			
30-50	Dense	8-15	Stiff	15-25				
>50	V. Dense	15-30	V. Stiff	>25				
		>30	Hard	and				

SUMMIT ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				SOIL BORING LOG			Boring #:	SE-SB-235		
Drilling Co: SUMMIT Geoengineering Services				Project: MAINE ENERGY PHASE II SUPPORT			Project #:	12-3259.1		
Personnel: Craig Coolidge				Location: BIDDEFORD, ME			Sheet:	1 of 1		
Summit Staff: John Cressey							Chkd by:	JKC		
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH						
Vehicle:	AMS	Type:	Concrete punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model:	9500-VTR	Hammer:	140LR	See Report						
Method:	H.S.A.	Fall:	24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	Blows/6 in.	SAMPLE DESCRIPTION		Stratum	Field Screening (ppmv)		
2	SS-1	30/29	0-2.5'		0-3" LOAM 3-10" SAND & GRAVEL (Fill) 10-15" CONCRETE & BRICK 15-29" Brown SAND & GRAVEL		LOAM			
							SAND & GRAVEL			
							DEMO DEBRIS			
	SS-2	30/30	2.5-5'		0-30" Reddish-brown SAND & GRAVEL		SAND & GRAVEL			
4	SS-3	24/22	5-7'		0-22" Reddish-brown SAND & GRAVEL					
	SS-4	24/20	7-9'		0-20" Reddish-brown SAND & GRAVEL					
	SS-5	24/20	9-11'		0-20" Reddish-brown SAND & GRAVEL					
6										
					CEASE @ 11'					
8										
10										
12										
14										
16										
18										
20										
Granular Soils		Cohesive Soils		% Composition		NOTES:				
Blows/ft.	Density	Blows/ft.	Consistency							
0-4	V. Loose	<2	V. soft	<5%	trace	1. Cease @ 11'				
4-10	Loose	2-4	Soft	5-15	little	2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')				
10-30	Compact	4-8	Firm	15-25	some	3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)				
30-50	Dense	8-15	Stiff	>25	and					
>50	V. Dense	15-30	V. Stiff							
		>30	Hard							

SUMMIT				SOIL BORING LOG			Boring #:	SE-SB-236		
ENVIRONMENTAL CONSULTANTS, INC. 640 Main Street Lewiston, Maine 04240				Project:	MAINE ENERGY PHASE II SUPPORT		Project #:	12-3259.1		
				Location:	BIDDEFORD, ME		Sheet:	1 of 1		
							Chkd by:	JKC		
Drilling Co: SUMMIT Geoengineering Services				Boring Location: see location map						
Personnel: Craig Coolidge				Top of PVC Casing Elevation:						
Summit Staff: John Cressey				Date started: 2/20/2013 Date Completed: 2/20/2013						
DRILLING METHOD		SAMPLER		ESTIMATED GROUND WATER DEPTH						
Vehicle:	AMS	Type:	Concrete punc	Date	Depth (ft)	Reference	Groundwater Elevation (ft)			
Model:	9500-VTR	Hammer:	140LR	See Report						
Method:	H.S.A.	Fall:	24"							
Depth (ft.)	No.	Pen/Rec (in)	Depth (ft)	SAMPLE DESCRIPTION			Stratum	Field Screening (ppmv)		
				Blows/6 in.						
	SS-1	24/16	0-2'		0-3" ASPHALT		ASPHALT			
					3-16" Brown SAND & GRAVEL (Fill)					
2							SAND & GRAVEL			
	SS-2	24/20	2-4'		0-12" Brown SAND & GRAVEL (Fill)					
					12-15 Brown SAND & GRAVEL w/ Brick		SAND, GRAVEL, DEBRIS			
4					15-20" Brown SAND & GRAVEL					
	SS-3	24/20	4-6'		0-8" Brown SAND & GRAVEL		SAND & GRAVEL			
					8-20" Brown SAND & ASH w/ Clinker					
6										
	SS-4	24/20	6-8'		0-20" Brown SAND & ASH w/ Clinker & Concrete					
							SAND, GRAVEL, DEBRIS			
8										
	SS-5	24/18	8-10'		0-18" Brown SAND & ASH w/ Clinker & Concrete					
10										
	SS-6	24/16	10-12'		0-6" Brown SAND & ASH w/ Clinker & Concrete					
					6-16" ROCK FRAGMENTS		ROCK			
12										
	SS-7	24/18	12-14'		0-18" Brown GRAVEL & ROCK, wet		GRAVEL & ROCK			
14					Cease @ 14'					
16										
18										
20										
Granular Soils		Cohesive Soils		% Composition						
Blows/ft.	Density	Blows/ft.	Consistency	NOTES:						
0-4	V. Loose	<2	V. soft	<5%	trace					
4-10	Loose	2-4	Soft	5-15	little					
10-30	Compact	4-8	Firm	15-25	some					
30-50	Dense	8-15	Stiff	>25	and					
>50	V. Dense	15-30	V. Stiff							
		>30	Hard							

1. Cease @ 14'
2. Samples taken at the following intervals: (0-2')(2-4')(4-6')(6-8')(8-10')(10-12')(12-14')
3. Samples to be analyzed for PCB's using the Soxhlet Extraction method (EPA 3540C)

Appendix C

Analytical Results

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: 12001150

Job ID: 25123

Date Received: 10/4/12

Project: MERC 12001150

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

A handwritten signature in black ink that appears to read "Cliff Chase".

Cliff Chase
Partner, Technical Director

Date of Approval: 10/23/2012
Total number of pages: 136

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SB-1 2.5-5	Solid	10/2/2012 15:30	25123-001	Solid Digestion for ICP Analysis Lead in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VPH in solids by MA DEP Method
CA-SB-2 0-1	Solid	10/2/2012 15:00	25123-002	EPH in solids by MADEP Method Solid Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-3 10-12.5	Solid	10/2/2012 13:10	25123-003	PCBs in soil by 8082 EPH in solids by MADEP Method Solid Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-4 0-2	Solid	10/2/2012 11:50	25123-004	Acid & Base/Neutral Extractables in solid by 8270 EPH in solids by MADEP Method Ranges only Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SB-4 0-2	Solid	10/2/2012 11:50	25123-004	Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260 VPH Ranges in solids by MA DEP Method
CA-SB-5 5-10	Solid	10/3/2012 13:30	25123-005	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260
CA-SB-6 0-5	Solid	10/3/2012 9:00	25123-006	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-6 12.5-15	Solid	10/3/2012 8:50	25123-007	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 EPH in solids by MADEP Method Ranges only Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SB-6 12.5-15	Solid	10/3/2012 8:50	25123-007	Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260 VPH Ranges in solids by MA DEP Method
CA-SB-7 7.5-10	Solid	10/2/2012 16:15	25123-008	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 EPH in solids by MADEP Method Ranges only Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260 VPH Ranges in solids by MA DEP Method
CA-SB-8 0-2.5	Solid	10/3/2012 8:50	25123-009	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-8 5-10	Solid	10/3/2012 10:00	25123-010	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SB-8 5-10	Solid	10/3/2012 10:00	25123-010	Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-9 0-4	Solid	10/3/2012 12:10	25123-011	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-10 10-15	Solid	10/3/2012 12:30	25123-012	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260
CA-SB-11 10-15	Solid	10/3/2012 8:30	25123-013	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 EPH in solids by MADEP Method Ranges only Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260 VPH Ranges in solids by MA DEP Method
SB-DUP-2	Solid	10/3/2012 0:00	25123-014	Acid & Base/Neutral Extractables in solid by 8270

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
SB-DUP-2	Solid	10/3/2012 0:00	25123-014	EPH in solids by MADEP Method Ranges only Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260 VPH Ranges in solids by MA DEP Method
CA-SB-12 15-18	Solid	10/2/2012 8:45	25123-015	PCBs in soil by 8082 EPH in solids by MADEP Method Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-13 10-12	Solid	10/2/2012 9:10	25123-016	PCBs in soil by 8082 EPH in solids by MADEP Method Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
SB-DUP-1	Solid	10/2/2012 0:00	25123-017	PCBs in soil by 8082 EPH in solids by MADEP Method Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Barium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Mercury in solids by 7471 Lead in solids by 6010 Selenium in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-14 0-3	Solid	10/2/2012 10:15	25123-018	EPH in solids by MADEP Method Percent Dry Matter for Sample Calc by SM2540B,G VPH in solids by MA DEP Method
CA-SB-15 5-5.5	Solid	10/2/2012 10:00	25123-019	EPH in solids by MADEP Method Percent Dry Matter for Sample Calc by SM2540B,G VPH in solids by MA DEP Method

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SS-1	Solid	10/3/2012 11:20	25123-020	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-2	Solid	10/3/2012 12:10	25123-021	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-3	Solid	10/3/2012 12:15	25123-022	PCBs in soil by 8082 PAHs in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SS-3	Solid	10/3/2012 12:15	25123-022	Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-4	Solid	10/3/2012 12:45	25123-023	PCBs in soil by 8082 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-5	Solid	10/3/2012 13:00	25123-024	PCBs in soil by 8082 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-CC-1	Solid	10/3/2012 9:30	25123-025	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-CC-1	Solid	10/3/2012 9:30	25123-025	Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260
CA-CC-2	Solid	10/3/2012 9:58	25123-026	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260
CA-CC-3	Solid	10/3/2012 10:40	25123-027	PCBs in soil by 8082 Acid & Base/Neutral Extractables in solid by 8270 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G VOCs in solids by 8260
CA-CC-4	Solid	10/3/2012 11:25	25123-028	PCBs in soil by 8082 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-CC-4	Solid	10/3/2012 11:25	25123-028	Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
CA-CC-5	Solid	10/3/2012 11:48	25123-029	PCBs in soil by 8082 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
STACK	Solid	10/3/2012 14:50	25123-030	PCBs in soil by 8082 Soil Digestion for ICP Analysis Silver in solids by 6010 Arsenic in solids by 6010 Beryllium in solids by 6010 Cadmium in solids by 6010 Chromium in solids by 6010 Copper in solids by 6010 Mercury in solids by 7471 Nickel in solids by 6010 Lead in solids by 6010 Antimony in solids by 6010 Selenium in solids by 6010 Thallium in solids by 6010 Zinc in solids by 6010 Percent Dry Matter for Sample Calc by SM2540B,G
Trip Blank	Solid	10/3/2012 0:00	25123-031	VOCs in solids by 8260
CA-SB-2 0-1 Re-Extraction	Solid	10/2/2012 15:00	25123-032	EPH in solids by MADEP Method Percent Dry Matter for Sample Calc by SM2540B,G

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	11:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
bromomethane				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
diethyl ether				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
acetone				< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
methylene chloride				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
2-butanone (MEK)				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
tetrahydrofuran (THF)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
2-hexanone				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B
m&p-xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 1:46	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Sampled: 10/2/12 11:50

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
Surrogate Recovery										
Limits										
dibromofluoromethane SUR	102	78-114	%	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
toluene-D8 SUR	99	88-110	%	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
4-bromofluorobenzene SUR	103	86-115	%	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B
a,a,a-trifluorotoluene SUR	99	70-130	%	1	LMM	10/8/12	5548	10/11/12	1:46	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	13:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
bromomethane				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
diethyl ether				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
acetone				< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
methylene chloride				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
2-butanone (MEK)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
tetrahydrofuran (THF)				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
2-hexanone				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B
m&p-xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:18	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Sampled: 10/3/12 13:30

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2,4-trichlorobenzene	0.2	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
Surrogate Recovery										
dibromofluoromethane SUR	102	78-114	%	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
toluene-D8 SUR	101	88-110	%	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
4-bromofluorobenzene SUR	99	86-115	%	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
a,a,a-trifluorotoluene SUR	100	70-130	%	1	LMM	10/8/12	5548	10/11/12	2:18	SW5035A8260B
Limits										

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	8:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
bromomethane				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
diethyl ether				< 0.6	0.6	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
acetone				< 3	3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
methylene chloride				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
2-butanone (MEK)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
tetrahydrofuran (THF)				< 0.6	0.6	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
2-hexanone				< 0.6	0.6	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
m&p xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	8:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
o-xylene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
styrene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
bromoform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
isopropylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,1,2,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2,3-trichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
n-propylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
bromobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,3,5-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
2-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
4-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
tert-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2,4-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
sec-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,3-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
4-isopropyltoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,4-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
n-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2,4-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
hexachlorobutadiene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
naphthalene				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
1,2,3-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
Surrogate Recovery												
						Limits						
dibromofluoromethane SUR				103	78-114	%	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
toluene-D8 SUR				101	88-110	%	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
4-bromofluorobenzene SUR				104	86-115	%	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B
a,a,a-trifluorotoluene SUR				97	70-130	%	1	LMM	10/8/12	5548	10/11/12 2:49	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	16:15	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
bromomethane				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
diethyl ether				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
acetone				< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
methylene chloride				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
2-butanone (MEK)				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
tetrahydrofuran (THF)				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
2-hexanone				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B
m&p xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	16:15	Reporting	Instr	Dil'n	Prep	Analysis			Reference
				Result	Limit	Units	Factor	Analyst	Date	Date	Time
o-xylene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
styrene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
bromoform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
isopropylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,1,2,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2,3-trichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
n-propylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
bromobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,3,5-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
2-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
4-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
tert-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2,4-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
sec-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,3-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
4-isopropyltoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,4-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
n-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2-dibromo-3-chloropropane (DBCP)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2,4-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
hexachlorobutadiene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
naphthalene				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
1,2,3-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:21
Surrogate Recovery											
						Limits					
dibromofluoromethane SUR				100	78-114	%	1	LMM	10/8/12	5548	10/11/12 3:21
toluene-D8 SUR				101	88-110	%	1	LMM	10/8/12	5548	10/11/12 3:21
4-bromofluorobenzene SUR				105	86-115	%	1	LMM	10/8/12	5548	10/11/12 3:21
a,a,a-trifluorotoluene SUR				98	70-130	%	1	LMM	10/8/12	5548	10/11/12 3:21

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	12:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
bromomethane				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
diethyl ether				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
acetone				< 3	3	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
methylene chloride				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
2-butanone (MEK)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
tetrahydrofuran (THF)				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
2-hexanone				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B
m&p-xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 3:52	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:30		Reporting		Instr Dil'n		Prep		Analysis			
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
o-xylene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
styrene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
bromoform		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
isopropylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,1,2,2-tetrachloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2,3-trichloropropane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
n-propylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
bromobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,3,5-trimethylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
2-chlorotoluene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
4-chlorotoluene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
tert-butylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2,4-trimethylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
sec-butylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,3-dichlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
4-isopropyltoluene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,4-dichlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2-dichlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
n-butylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2,4-trichlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
hexachlorobutadiene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
naphthalene		< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
1,2,3-trichlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
Surrogate Recovery		Limits									
dibromofluoromethane SUR		100	78-114	%	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
toluene-D8 SUR		96	88-110	%	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
4-bromofluorobenzene SUR		103	86-115	%	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B
a,a,a-trifluorotoluene SUR		101	70-130	%	1	LMM	10/8/12	5548	10/11/12	3:52	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	8:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	
dichlorodifluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
chloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
vinyl chloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
bromomethane				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
chloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
trichlorofluoromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
diethyl ether				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
acetone				< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
methylene chloride				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
carbon disulfide				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
methyl t-butyl ether (MTBE)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
trans-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
2-butanone (MEK)				< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
2,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
cis-1,2-dichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
chloroform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
bromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
tetrahydrofuran (THF)				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1,1-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
carbon tetrachloride				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,2-dichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
benzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
trichloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,2-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
bromodichloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
dibromomethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
4-methyl-2-pentanone (MIBK)				< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
cis-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
toluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
trans-1,3-dichloropropene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
2-hexanone				< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1,2-trichloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,3-dichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
tetrachloroethene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
dibromochloromethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,2-dibromoethane (EDB)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
chlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
1,1,1,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
ethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B
m&p-xylenes				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12 5:27	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:30

Parameter	Reporting		Instr	Dil'n	Prep	Analysis			Reference	
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
Surrogate Recovery										
Limits										
dibromofluoromethane SUR	101	78-114	%	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
toluene-D8 SUR	100	88-110	%	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
4-bromofluorobenzene SUR	104	86-115	%	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B
a,a,a-trifluorotoluene SUR	106	70-130	%	1	LMM	10/8/12	5548	10/11/12	5:27	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 0:00	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Date	Time		
dichlorodifluoromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
chloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
vinyl chloride			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
bromomethane			< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
chloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
trichlorofluoromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
diethyl ether			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
acetone			< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
methylene chloride			< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
carbon disulfide			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
methyl t-butyl ether (MTBE)			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
trans-1,2-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1-dichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
2-butanone (MEK)			< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
2,2-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
cis-1,2-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
chloroform			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
bromochloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
tetrahydrofuran (THF)			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1,1-trichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
carbon tetrachloride			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2-dichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
benzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
trichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
bromodichloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
dibromomethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
4-methyl-2-pentanone (MIBK)			< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
cis-1,3-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
toluene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
trans-1,3-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
2-hexanone			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1,2-trichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,3-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
tetrachloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
dibromochloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2-dibromoethane (EDB)			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
chlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1,1,2-tetrachloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
ethylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
m&p-xylenes			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 10/3/12 0:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
Surrogate Recovery		Limits								
dibromofluoromethane SUR	101	78-114	%	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
toluene-D8 SUR	103	88-110	%	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
4-bromofluorobenzene SUR	100	86-115	%	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B
a,a,a-trifluorotoluene SUR	99	70-130	%	1	LMM	10/8/12	5548	10/11/12	4:55	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 9:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Date	Time		
dichlorodifluoromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
chloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
vinyl chloride			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
bromomethane			< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
chloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
trichlorofluoromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
diethyl ether			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
acetone			< 2	2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
methylene chloride			< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
carbon disulfide			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
methyl t-butyl ether (MTBE)			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
trans-1,2-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1-dichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
2-butanone (MEK)			< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
2,2-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
cis-1,2-dichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
chloroform			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
bromochloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
tetrahydrofuran (THF)			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1,1-trichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
carbon tetrachloride			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2-dichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
benzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
trichloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
bromodichloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
dibromomethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
4-methyl-2-pentanone (MIBK)			< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
cis-1,3-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
toluene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
trans-1,3-dichloropropene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
2-hexanone			< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1,2-trichloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,3-dichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
tetrachloroethene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
dibromochloromethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2-dibromoethane (EDB)			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
chlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1,1,2-tetrachloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
ethylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
m&p-xylenes			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 9:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Date	Time		
o-xylene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
styrene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
bromoform			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
isopropylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,1,2,2-tetrachloroethane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2,3-trichloropropane			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
n-propylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
bromobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,3,5-trimethylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
2-chlorotoluene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
4-chlorotoluene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
tert-butylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2,4-trimethylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
sec-butylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,3-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
4-isopropyltoluene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,4-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
n-butylbenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2,4-trichlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
hexachlorobutadiene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
naphthalene			0.4	0.2	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
1,2,3-trichlorobenzene			< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
Surrogate Recovery												
			Limits									
dibromofluoromethane SUR			101	78-114	%	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
toluene-D8 SUR			101	88-110	%	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
4-bromofluorobenzene SUR			106	86-115	%	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B
a,a,a-trifluorotoluene SUR			103	70-130	%	1	LMM	10/8/12	5548	10/11/12	4:24	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:58	Reporting	Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Time	Reference
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
bromomethane	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
diethyl ether	< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
acetone	< 2	2	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
methylene chloride	< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12 17:37 SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 9:58	Reporting	Instr	Dil'n	Prep	Analysis				
			Result	Limit	Units	Analyst	Date	Batch	Date	Time	Reference
o-xylene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
styrene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
bromoform			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
isopropylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,1,2,2-tetrachloroethane			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2,3-trichloropropane			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
n-propylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
bromobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,3,5-trimethylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
2-chlorotoluene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
4-chlorotoluene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
tert-butylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2,4-trimethylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
sec-butylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,3-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
4-isopropyltoluene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,4-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2-dichlorobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
n-butylbenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2,4-trichlorobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
hexachlorobutadiene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
naphthalene			< 0.2	0.2	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
1,2,3-trichlorobenzene			< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
Surrogate Recovery											
dibromofluoromethane SUR			99	78-114	%	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
toluene-D8 SUR			102	88-110	%	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
4-bromofluorobenzene SUR			107	86-115	%	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
a,a,a-trifluorotoluene SUR			101	70-130	%	1	LMM 10/8/12	5548	10/12/12	17:37	SW5035A8260B
Limits											

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Sampled:	10/3/12 10:40	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
dichlorodifluoromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
chloromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
vinyl chloride		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
bromomethane		< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
chloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
trichlorofluoromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
diethyl ether		< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
acetone		< 2	2	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1-dichloroethene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
methylene chloride		< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
carbon disulfide		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
methyl t-butyl ether (MTBE)		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
trans-1,2-dichloroethene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1-dichloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
2-butanone (MEK)		< 0.3	0.3	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
2,2-dichloropropane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
cis-1,2-dichloroethene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
chloroform		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
bromochloromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
tetrahydrofuran (THF)		< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1,1-trichloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1-dichloropropene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
carbon tetrachloride		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2-dichloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
benzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
trichloroethene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2-dichloropropane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
bromodichloromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
dibromomethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
4-methyl-2-pentanone (MIBK)		< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
cis-1,3-dichloropropene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
toluene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
trans-1,3-dichloropropene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
2-hexanone		< 0.4	0.4	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1,2-trichloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,3-dichloropropane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
tetrachloroethene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
dibromochloromethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2-dibromoethane (EDB)		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
chlorobenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1,1,2-tetrachloroethane		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
ethylbenzene		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
m&p-xylenes		< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	10:40	Reporting	Instr	Dil'n	Prep	Analysis					
				Result	Limit	Units	Factor	Analyst	Date	Date	Time	Reference	
o-xylene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
styrene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
bromoform				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
isopropylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,1,2,2-tetrachloroethane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2,3-trichloropropane				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
n-propylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
bromobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,3,5-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
2-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
4-chlorotoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
tert-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2,4-trimethylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
sec-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,3-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
4-isopropyltoluene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,4-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2-dichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
n-butylbenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2,4-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
hexachlorobutadiene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
naphthalene				< 0.2	0.2	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
1,2,3-trichlorobenzene				< 0.1	0.1	ug/g	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
Surrogate Recovery													
dibromofluoromethane SUR				96	78-114	%	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
toluene-D8 SUR				103	88-110	%	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
4-bromofluorobenzene SUR				102	86-115	%	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
a,a,a-trifluorotoluene SUR				100	70-130	%	1	LMM	10/8/12	5548	10/12/12	18:08	SW5035A8260B
Limits													

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-031

Sample ID: Trip Blank

Matrix: Solid

Sampled: 10/3/12 0:00

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
dichlorodifluoromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
chloromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
vinyl chloride	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
bromomethane	< 0.2	0.2	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
chloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
trichlorofluoromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
diethyl ether	< 0.5	0.5	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
acetone	< 2	2	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1-dichloroethene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
methylene chloride	< 0.2	0.2	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
carbon disulfide	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
trans-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1-dichloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
2-butanone (MEK)	< 0.3	0.3	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
2,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
cis-1,2-dichloroethene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
chloroform	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
bromochloromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
tetrahydrofuran (THF)	< 0.5	0.5	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1,1-trichloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1-dichloropropene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
carbon tetrachloride	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2-dichloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
benzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
trichloroethene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2-dichloropropane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
bromodichloromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
dibromomethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
4-methyl-2-pentanone (MIBK)	< 0.4	0.4	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
cis-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
toluene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
trans-1,3-dichloropropene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
2-hexanone	< 0.5	0.5	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1,2-trichloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,3-dichloropropane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
tetrachloroethene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
dibromochloromethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2-dibromoethane (EDB)	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
chlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1,1,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
ethylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-031

Sample ID: Trip Blank

Matrix: Solid

Sampled: 10/3/12 0:00

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
o-xylene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
styrene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
bromoform	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
isopropylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,1,2,2-tetrachloroethane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2,3-trichloropropane	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
n-propylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
bromobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,3,5-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
2-chlorotoluene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
4-chlorotoluene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
tert-butylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2,4-trimethylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
sec-butylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,3-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
4-isopropyltoluene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,4-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2-dichlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
n-butylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2-dibromo-3-chloropropane (DBCP)	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2,4-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
hexachlorobutadiene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
naphthalene	< 0.2	0.2	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
1,2,3-trichlorobenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
Surrogate Recovery									
dibromofluoromethane SUR	104	78-114	%	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
toluene-D8 SUR	101	88-110	%	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
4-bromofluorobenzene SUR	99	86-115	%	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
a,a,a-trifluorotoluene SUR	95	70-130	%	1	LMM 10/8/12	5548	10/10/12	23:40	SW5035A8260B
Limits									

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-001

Sample ID: CA-SB-1 2.5-5

Matrix: Solid Percent Dry: 84.1% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.62 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/2/12 15:30

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 5	5	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
Unadjusted C9-C12 Aliphatics	24	5	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
benzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
toluene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
o-xylene	< 0.1	0.1	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
naphthalene	< 0.2	0.2	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
C5-C8 Aliphatics	< 5	5	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
C9-C12 Aliphatics	23	5	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
C9-C10 Aromatics	< 5	5	ug/g	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
Surrogate Recovery									
Limits									
2,5-dibromotoluene as Aromatic SUR	104	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
2,5-dibromotoluene as Aliphatic SUR	100	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH
a,a,a-trifluorotoluene SUR	115	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:02	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid Percent Dry: 97.2% Results expressed on a dry weight basis.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/2/12 11:50

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 3	3	ug/g	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH
Unadjusted C9-C12 Aliphatics	< 3	3	ug/g	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH
C9-C10 Aromatics	< 3	3	ug/g	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH
Surrogate Recovery									
Limits									
2,5-dibromotoluene as Aromatic SUR	101	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH
2,5-dibromotoluene as Aliphatic SUR	103	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH
a,a,a-trifluorotoluene SUR	100	70-130	%	1	LMM 10/8/12	5549	10/9/12	21:35	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid Percent Dry: 77% Results expressed on a dry weight basis.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/3/12 8:50

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 6	6	ug/g	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH
Unadjusted C9-C12 Aliphatics	< 6	6	ug/g	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH
C9-C10 Aromatics	< 6	6	ug/g	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH
Surrogate Recovery										
Limits										
2,5-dibromotoluene as Aromatic SUR	102	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH
2,5-dibromotoluene as Aliphatic SUR	99	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH
a,a,a-trifluorotoluene SUR	97	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:08	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid Percent Dry: 88.4% Results expressed on a dry weight basis.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/2/12 16:15

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH
Unadjusted C9-C12 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH
C9-C10 Aromatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH
Surrogate Recovery										
Limits										
2,5-dibromotoluene as Aromatic SUR	100	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH
2,5-dibromotoluene as Aliphatic SUR	102	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH
a,a,a-trifluorotoluene SUR	100	70-130	%	1	LMM	10/8/12	5549	10/9/12	22:41	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid Percent Dry: 78.5% Results expressed on a dry weight basis.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/3/12 8:30

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH
Unadjusted C9-C12 Aliphatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH
C9-C10 Aromatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH
Surrogate Recovery										
Limits										
2,5-dibromotoluene as Aromatic SUR	102	70-130	%	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH
2,5-dibromotoluene as Aliphatic SUR	104	70-130	%	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH
a,a,a-trifluorotoluene SUR	116	70-130	%	1	LMM	10/8/12	5549	10/10/12	1:58	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid Percent Dry: 77.6% Results expressed on a dry weight basis.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/3/12 0:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH
Unadjusted C9-C12 Aliphatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH
C9-C10 Aromatics	< 5	5	ug/g	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH
Surrogate Recovery										
Limits										
2,5-dibromotoluene as Aromatic SUR	101	70-130	%	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH
2,5-dibromotoluene as Aliphatic SUR	105	70-130	%	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH
a,a,a-trifluorotoluene SUR	111	70-130	%	1	LMM	10/8/12	5549	10/10/12	0:52	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-018

Sample ID: CA-SB-14 0-3

Matrix: Solid Percent Dry: 94.8% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.69 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/2/12 10:15

Parameter	Reporting		Instr Dil'n		Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
Unadjusted C9-C12 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
C5-C8 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
C9-C12 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
C9-C10 Aromatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
Surrogate Recovery										
2,5-dibromotoluene as Aromatic SUR	99	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
2,5-dibromotoluene as Aliphatic SUR	102	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH
a,a,a-trifluorotoluene SUR	103	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:47	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-019

Sample ID: CA-SB-15 5-5.5

Matrix: Solid Percent Dry: 88.1% Results expressed on a dry weight basis.

Samples prepared in methanol at a ratio of 0.64 mL MeOH/g soil.

Received on ice at 0°C, in satisfactory condition.

Sampled: 10/2/12 10:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
Unadjusted C9-C12 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
methyl t-butyl ether (MTBE)	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
benzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
toluene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
ethylbenzene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
m&p-xylenes	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
o-xylene	< 0.1	0.1	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
naphthalene	< 0.2	0.2	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
C5-C8 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
C9-C12 Aliphatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
C9-C10 Aromatics	< 4	4	ug/g	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
Surrogate Recovery										
Limits										
2,5-dibromotoluene as Aromatic SUR	100	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
2,5-dibromotoluene as Aliphatic SUR	107	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH
a,a,a-trifluorotoluene SUR	101	70-130	%	1	LMM	10/8/12	5549	10/9/12	23:14	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	11:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
aniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
phenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-chlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
bis(2-chloroethyl)ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
1,3-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
1,4-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
1,2-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzyl alcohol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
bis(2-chloroisopropyl) ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
hexachloroethane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
N-nitroso-di-N-propylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
nitrobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
isophorone				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-nitrophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4-dimethylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
bis(2-chloroethoxy)methane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4-dichlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
1,2,4-trichlorobenzene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
naphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzoic acid				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-chloroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
hexachlorobutadiene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-chloro-3-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-methylnaphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
hexachlorocyclopentadiene				< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4,6-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4,5-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-chloronaphthalene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
acenaphthylene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
dimethylphthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,6-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
acenaphthene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
3-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4-dinitrophenol				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
dibenzofuran				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-nitrophenol				< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
fluorene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
diethyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	11:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
4-chlorophenyl phenyl ether				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-nitroaniline				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4,6-dinitro-2-methylphenol				< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
azobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
N-nitrosodiphenylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
4-bromophenyl phenyl ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
hexachlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
pentachlorophenol				< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
phenanthrene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
anthracene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
carbazole				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
di-n-butylphthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
fluoranthene				0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzidine				< 3	3	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
pyrene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
butyl benzyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzo(a)anthracene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
chrysene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
3,3'-dichlorobenzidine				< 3	3	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
bis(2-ethylhexyl)phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
di-n-octyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzo(b)fluoranthene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzo(k)fluoranthene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzo(a)pyrene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
indeno(1,2,3-cd)pyrene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
dibenzo(a,h)anthracene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
benzo(g,h,i)perylene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
Surrogate Recovery												
						Limits						
2-fluorophenol SUR				52	21-100	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
phenol-D5 SUR				52	10-102	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2,4,6-tribromophenol SUR				70	10-123	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
nitrobenzene-D5 SUR				52	35-114	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
2-fluorobiphenyl SUR				68	43-116	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D
p-terphenyl-D14 SUR				69	33-141	%	1	AJD 10/10/12	5556	10/12/12	15:33	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	13:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
aniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
phenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-chlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
bis(2-chloroethyl)ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
1,3-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
1,4-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
1,2-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzyl alcohol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
bis(2-chloroisopropyl) ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
hexachloroethane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
N-nitroso-di-N-propylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
nitrobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
isophorone				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-nitrophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4-dimethylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
bis(2-chloroethoxy)methane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4-dichlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
1,2,4-trichlorobenzene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
naphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzoic acid				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-chloroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
hexachlorobutadiene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-chloro-3-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-methylnaphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
hexachlorocyclopentadiene				< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4,6-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4,5-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-chloronaphthalene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
acenaphthylene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
dimethylphthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,6-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
acenaphthene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
3-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4-dinitrophenol				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
dibenzofuran				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-nitrophenol				< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
fluorene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
diethyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	13:30	Reporting	Instr	Dil'n	Prep	Analysis				
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-nitroaniline				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4,6-dinitro-2-methylphenol				< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
azobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
N-nitrosodiphenylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
4-bromophenyl phenyl ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
hexachlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
pentachlorophenol				< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
phenanthrene				0.09	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
anthracene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
carbazole				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
di-n-butylphthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
fluoranthene				0.21	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzidine				< 3	3	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
pyrene				0.19	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
butyl benzyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzo(a)anthracene				0.12	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
chrysene				0.13	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
3,3'-dichlorobenzidine				< 3	3	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
bis(2-ethylhexyl)phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
di-n-octyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzo(b)fluoranthene				0.14	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzo(k)fluoranthene				0.15	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzo(a)pyrene				0.12	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
indeno(1,2,3-cd)pyrene				0.06	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
dibenzo(a,h)anthracene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
benzo(g,h,i)perylene				0.06	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
Surrogate Recovery												
						Limits						
2-fluorophenol SUR				47	21-100	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
phenol-D5 SUR				47	10-102	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2,4,6-tribromophenol SUR				53	10-123	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
nitrobenzene-D5 SUR				47	35-114	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
2-fluorobiphenyl SUR				60	43-116	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D
p-terphenyl-D14 SUR				59	33-141	%	1	AJD 10/10/12	5556	10/12/12	21:45	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-006

Sample ID: CA-SB-6 0-5

Matrix: Solid

Percent Dry: 93.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:00		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
naphthalene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
2-methylnaphthalene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
acenaphthylene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
acenaphthene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
dibenzofuran	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
fluorene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
phenanthrene	0.6	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
anthracene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
fluoranthene	1.2	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
pyrene	1.0	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
benzo(a)anthracene	0.6	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
chrysene	0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
benzo(b)fluoranthene	0.6	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
benzo(k)fluoranthene	0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
benzo(a)pyrene	0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
dibenzo(a,h)anthracene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
benzo(g,h,i)perylene	< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
Surrogate Recovery											
2-fluorobiphenyl SUR	93	43-116	%	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
o-terphenyl SUR	96	33-141	%	1	AJD	10/10/12	5555	10/11/12	18:56	SW3550B8270D	
Limits											

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 8:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
aniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
phenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-chlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
bis(2-chloroethyl)ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
1,3-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
1,4-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
1,2-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzyl alcohol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
bis(2-chloroisopropyl) ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
hexachloroethane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
N-nitroso-di-N-propylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
nitrobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
isophorone			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-nitrophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4-dimethylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
bis(2-chloroethoxy)methane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4-dichlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
1,2,4-trichlorobenzene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
naphthalene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzoic acid			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-chloroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
hexachlorobutadiene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-chloro-3-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-methylnaphthalene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
hexachlorocyclopentadiene			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4,6-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4,5-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-chloronaphthalene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
acenaphthylene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
dimethylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,6-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
acenaphthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
3-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4-dinitrophenol			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
dibenzofuran			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-nitrophenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
fluorene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
diethyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 8:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
4-chlorophenyl phenyl ether			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-nitroaniline			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4,6-dinitro-2-methylphenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
azobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
N-nitrosodiphenylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
4-bromophenyl phenyl ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
hexachlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
pentachlorophenol			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
phenanthrene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
anthracene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
carbazole			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
di-n-butylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
fluoranthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzidine			< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
pyrene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
butyl benzyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzo(a)anthracene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
chrysene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
3,3'-dichlorobenzidine			< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
bis(2-ethylhexyl)phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
di-n-octyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzo(b)fluoranthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzo(k)fluoranthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzo(a)pyrene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
indeno(1,2,3-cd)pyrene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
dibenzo(a,h)anthracene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
benzo(g,h,i)perylene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
Surrogate Recovery												
			Limits									
2-fluorophenol SUR			56	21-100	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
phenol-D5 SUR			53	10-102	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2,4,6-tribromophenol SUR			61	10-123	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
nitrobenzene-D5 SUR			53	35-114	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
2-fluorobiphenyl SUR			64	43-116	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D
p-terphenyl-D14 SUR			58	33-141	%	1	AJD	10/10/12	5556	10/12/12	16:47	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Sampled:	10/2/12 16:15	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
aniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
phenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-chlorophenol		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
bis(2-chloroethyl)ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
1,3-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
1,4-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
1,2-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzyl alcohol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
bis(2-chloroisopropyl) ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
hexachloroethane		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
N-nitroso-di-N-propylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
nitrobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
isophorone		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-nitrophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4-dimethylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
bis(2-chloroethoxy)methane		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4-dichlorophenol		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
1,2,4-trichlorobenzene		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
naphthalene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzoic acid		< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-chloroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
hexachlorobutadiene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-chloro-3-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-methylnaphthalene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
hexachlorocyclopentadiene		< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4,6-trichlorophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4,5-trichlorophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-chloronaphthalene		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-nitroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
acenaphthylene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
dimethylphthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,6-dinitrotoluene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4-dinitrotoluene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
acenaphthene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
3-nitroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4-dinitrophenol		< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
dibenzofuran		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-nitrophenol		< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
fluorene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
diethyl phthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Sampled: 10/2/12 16:15	Reporting	Instr	Dil'n	Prep	Analysis					
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-nitroaniline	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4,6-dinitro-2-methylphenol	< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
azobenzene	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
N-nitrosodiphenylamine	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
4-bromophenyl phenyl ether	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
hexachlorobenzene	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
pentachlorophenol	< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
phenanthrene	0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
anthracene	< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
carbazole	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
di-n-butylphthalate	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
fluoranthene	0.08	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzidine	< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
pyrene	0.09	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
butyl benzyl phthalate	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzo(a)anthracene	0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
chrysene	0.07	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
3,3'-dichlorobenzidine	< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
bis(2-ethylhexyl)phthalate	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
di-n-octyl phthalate	< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzo(b)fluoranthene	< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzo(k)fluoranthene	0.09	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzo(a)pyrene	0.08	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
indeno(1,2,3-cd)pyrene	0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
dibenzo(a,h)anthracene	< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
benzo(g,h,i)perylene	0.07	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
Surrogate Recovery										
Limits										
2-fluorophenol SUR	55	21-100	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
phenol-D5 SUR	52	10-102	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2,4,6-tribromophenol SUR	63	10-123	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
nitrobenzene-D5 SUR	52	35-114	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
2-fluorobiphenyl SUR	64	43-116	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D
p-terphenyl-D14 SUR	65	33-141	%	1	AJD	10/10/12	5556	10/12/12	16:10	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-009

Sample ID: CA-SB-8 0-2.5

Matrix: Solid

Percent Dry: 97.7% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 8:50	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time
naphthalene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
2-methylnaphthalene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
acenaphthylene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
acenaphthene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
dibenzofuran			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
fluorene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
phenanthrene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
anthracene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
fluoranthene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
pyrene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
benzo(a)anthracene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
chrysene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
benzo(b)fluoranthene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
benzo(k)fluoranthene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
benzo(a)pyrene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
indeno(1,2,3-cd)pyrene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
dibenzo(a,h)anthracene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
benzo(g,h,i)perylene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5555	10/11/12	17:04
Surrogate Recovery											
2-fluorobiphenyl SUR			86	43-116	%	1	AJD	10/10/12	5555	10/11/12	17:04
o-terphenyl SUR			89	33-141	%	1	AJD	10/10/12	5555	10/11/12	17:04
Limits											

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-010

Sample ID: CA-SB-8 5-10

Matrix: Solid

Percent Dry: 83.8% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	10:00	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time
naphthalene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
2-methylnaphthalene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
acenaphthylene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
acenaphthene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
dibenzofuran				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
fluorene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
phenanthrene				0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
anthracene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
fluoranthene				0.7	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
pyrene				0.8	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
benzo(a)anthracene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
chrysene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
benzo(b)fluoranthene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
benzo(k)fluoranthene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
benzo(a)pyrene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
indeno(1,2,3-cd)pyrene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
dibenzo(a,h)anthracene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
benzo(g,h,i)perylene				< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	19:33
Surrogate Recovery												
2-fluorobiphenyl SUR				89	43-116	%	1	AJD	10/10/12	5555	10/11/12	19:33
o-terphenyl SUR				88	33-141	%	1	AJD	10/10/12	5555	10/11/12	19:33
Limits												

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:30

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
N-nitrosodimethylamine	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
aniline	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
phenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-chlorophenol	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
bis(2-chloroethyl)ether	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
1,3-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
1,4-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
1,2-dichlorobenzene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzyl alcohol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-methylphenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
bis(2-chloroisopropyl) ether	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
hexachloroethane	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
N-nitroso-di-N-propylamine	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-methylphenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
nitrobenzene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
isophorone	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-nitrophenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4-dimethylphenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
bis(2-chloroethoxy)methane	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4-dichlorophenol	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
1,2,4-trichlorobenzene	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
naphthalene	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzoic acid	< 6	6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-chloroaniline	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
hexachlorobutadiene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-chloro-3-methylphenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-methylnaphthalene	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
hexachlorocyclopentadiene	< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4,6-trichlorophenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4,5-trichlorophenol	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-chloronaphthalene	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-nitroaniline	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
acenaphthylene	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
dimethylphthalate	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,6-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4-dinitrotoluene	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
acenaphthene	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
3-nitroaniline	< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4-dinitrophenol	< 6	6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
dibenzofuran	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-nitrophenol	< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
fluorene	< 0.06	0.06	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D
diethyl phthalate	< 0.6	0.6	ug/g	1	AJD 10/10/12	5556	10/12/12	19:16	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Sampled:	10/3/12 12:30	Reporting	Instr	Dil'n	Prep	Analysis					
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-nitroaniline		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
4,6-dinitro-2-methylphenol		< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
azobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
N-nitrosodiphenylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
4-bromophenyl phenyl ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
hexachlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
pentachlorophenol		< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
phenanthrene		0.10	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
anthracene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
carbazole		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
di-n-butylphthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
fluoranthene		0.25	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzidine		< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
pyrene		0.28	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
butyl benzyl phthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzo(a)anthracene		0.21	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
chrysene		0.20	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
3,3'-dichlorobenzidine		< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
bis(2-ethylhexyl)phthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
di-n-octyl phthalate		< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzo(b)fluoranthene		0.18	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzo(k)fluoranthene		0.18	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzo(a)pyrene		0.20	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
indeno(1,2,3-cd)pyrene		0.11	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
dibenzo(a,h)anthracene		< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
benzo(g,h,i)perylene		0.10	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
Surrogate Recovery											
		Limits									
2-fluorophenol SUR		44	21-100	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
phenol-D5 SUR		43	10-102	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
2,4,6-tribromophenol SUR		59	10-123	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
nitrobenzene-D5 SUR		42	35-114	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
2-fluorobiphenyl SUR		51	43-116	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D
p-terphenyl-D14 SUR		58	33-141	%	1	AJD	10/10/12	5556	10/12/12	19:16	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 8:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time
N-nitrosodimethylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
aniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
phenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-chlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
bis(2-chloroethyl)ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
1,3-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
1,4-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
1,2-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
benzyl alcohol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
bis(2-chloroisopropyl) ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
hexachloroethane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
N-nitroso-di-N-propylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
4-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
nitrobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
isophorone			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-nitrophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4-dimethylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
bis(2-chloroethoxy)methane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4-dichlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
1,2,4-trichlorobenzene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
naphthalene			0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
benzoic acid			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
4-chloroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
hexachlorobutadiene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
4-chloro-3-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-methylnaphthalene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
hexachlorocyclopentadiene			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4,6-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4,5-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-chloronaphthalene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
acenaphthylene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
dimethylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,6-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
acenaphthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
3-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
2,4-dinitrophenol			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
dibenzofuran			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
4-nitrophenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
fluorene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54
diethyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 8:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
4-chlorophenyl phenyl ether			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
4-nitroaniline			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
4,6-dinitro-2-methylphenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
azobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
N-nitrosodiphenylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
4-bromophenyl phenyl ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
hexachlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
pentachlorophenol			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
phenanthrene			0.37	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
anthracene			0.08	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
carbazole			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
di-n-butylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
fluoranthene			0.38	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzidine			< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
pyrene			0.38	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
butyl benzyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzo(a)anthracene			0.20	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
chrysene			0.21	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
3,3'-dichlorobenzidine			< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
bis(2-ethylhexyl)phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
di-n-octyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzo(b)fluoranthene			0.13	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzo(k)fluoranthene			0.17	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzo(a)pyrene			0.17	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
indeno(1,2,3-cd)pyrene			0.07	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
dibenzo(a,h)anthracene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
benzo(g,h,i)perylene			0.08	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
Surrogate Recovery												
			Limits									
2-fluorophenol SUR			50	21-100	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
phenol-D5 SUR			46	10-102	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
2,4,6-tribromophenol SUR			68	10-123	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
nitrobenzene-D5 SUR			46	35-114	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
2-fluorobiphenyl SUR			60	43-116	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D
p-terphenyl-D14 SUR			58	33-141	%	1	AJD	10/10/12	5556	10/12/12	19:54	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 0:00	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
aniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
phenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-chlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
bis(2-chloroethyl)ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
1,3-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
1,4-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
1,2-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzyl alcohol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
bis(2-chloroisopropyl) ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
hexachloroethane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
N-nitroso-di-N-propylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
nitrobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
isophorone			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-nitrophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4-dimethylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
bis(2-chloroethoxy)methane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4-dichlorophenol			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
1,2,4-trichlorobenzene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
naphthalene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzoic acid			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-chloroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
hexachlorobutadiene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-chloro-3-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-methylnaphthalene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
hexachlorocyclopentadiene			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4,6-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4,5-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-chloronaphthalene			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
acenaphthylene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
dimethylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,6-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
acenaphthene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
3-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4-dinitrophenol			< 6	6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
dibenzofuran			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-nitrophenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
fluorene			< 0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
diethyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 0:00	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
4-chlorophenyl phenyl ether			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-nitroaniline			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4,6-dinitro-2-methylphenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
azobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
N-nitrosodiphenylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
4-bromophenyl phenyl ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
hexachlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
pentachlorophenol			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
phenanthrene			0.48	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
anthracene			0.11	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
carbazole			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
di-n-butylphthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
fluoranthene			0.64	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzidine			< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
pyrene			0.60	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
butyl benzyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzo(a)anthracene			0.33	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
chrysene			0.33	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
3,3'-dichlorobenzidine			< 4	4	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
bis(2-ethylhexyl)phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
di-n-octyl phthalate			< 0.6	0.6	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzo(b)fluoranthene			0.32	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzo(k)fluoranthene			0.23	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzo(a)pyrene			0.31	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
indeno(1,2,3-cd)pyrene			0.15	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
dibenzo(a,h)anthracene			0.06	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
benzo(g,h,i)perylene			0.15	0.06	ug/g	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
Surrogate Recovery												
			Limits									
2-fluorophenol SUR			50	21-100	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
phenol-D5 SUR			47	10-102	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2,4,6-tribromophenol SUR			66	10-123	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
nitrobenzene-D5 SUR			48	35-114	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
2-fluorobiphenyl SUR			62	43-116	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D
p-terphenyl-D14 SUR			57	33-141	%	1	AJD	10/10/12	5556	10/12/12	20:31	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-020

Sample ID: CA-SS-1

Matrix: Solid

Percent Dry: 82.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep	Analysis			Reference
	Result	Limit	Units	Factor		Date	Batch	Date	Time	
naphthalene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
2-methylnaphthalene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
acenaphthylene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
acenaphthene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
dibenzofuran	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
fluorene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
phenanthrene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
anthracene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
fluoranthene	0.8	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
pyrene	0.7	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
benzo(a)anthracene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
chrysene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
benzo(b)fluoranthene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
benzo(k)fluoranthene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
benzo(a)pyrene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
indeno(1,2,3-cd)pyrene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
dibenzo(a,h)anthracene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
benzo(g,h,i)perylene	< 0.6	0.6	ug/g	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
Surrogate Recovery										
2-fluorobiphenyl SUR	95	43-116	%	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
o-terphenyl SUR	100	33-141	%	1	AJD	10/10/12	5555	10/11/12	20:10	SW3550B8270D
Limits										

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-021

Sample ID: CA-SS-2

Matrix: Solid

Percent Dry: 91.8% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	12:10	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
naphthalene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
2-methylnaphthalene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
acenaphthylene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
acenaphthene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
dibenzofuran				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
fluorene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
phenanthrene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
anthracene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
fluoranthene				0.8	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
pyrene				0.8	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
benzo(a)anthracene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
chrysene				0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
benzo(b)fluoranthene				0.7	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
benzo(k)fluoranthene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
benzo(a)pyrene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
indeno(1,2,3-cd)pyrene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
dibenzo(a,h)anthracene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
benzo(g,h,i)perylene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
Surrogate Recovery												
2-fluorobiphenyl SUR				88	43-116	%	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
o-terphenyl SUR				90	33-141	%	1	AJD 10/10/12	5555	10/11/12	23:16	SW3550B8270D
Limits												

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-022

Sample ID: CA-SS-3

Matrix: Solid

Percent Dry: 93.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	12:15	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
naphthalene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
2-methylnaphthalene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
acenaphthylene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
acenaphthene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
dibenzofuran				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
fluorene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
phenanthrene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
anthracene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
fluoranthene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
pyrene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
benzo(a)anthracene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
chrysene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
benzo(b)fluoranthene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
benzo(k)fluoranthene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
benzo(a)pyrene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
indeno(1,2,3-cd)pyrene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
dibenzo(a,h)anthracene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
benzo(g,h,i)perylene				< 0.5	0.5	ug/g	AJD	10/10/12	5555	10/11/12	20:47	
Surrogate Recovery												
2-fluorobiphenyl SUR				90	43-116	%	1	AJD	10/10/12	5555	10/11/12	20:47
o-terphenyl SUR				92	33-141	%	1	AJD	10/10/12	5555	10/11/12	20:47
Limits												

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 9:30	Reporting	Instr	Dil'n	Prep	Analysis			Reference		
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
aniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
phenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-chlorophenol			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
bis(2-chloroethyl)ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
1,3-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
1,4-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
1,2-dichlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzyl alcohol			0.4	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
bis(2-chloroisopropyl) ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
hexachloroethane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
N-nitroso-di-N-propylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
nitrobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
isophorone			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-nitrophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4-dimethylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
bis(2-chloroethoxy)methane			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4-dichlorophenol			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
1,2,4-trichlorobenzene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
naphthalene			0.32	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzoic acid			< 5	5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-chloroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
hexachlorobutadiene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-chloro-3-methylphenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-methylnaphthalene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
hexachlorocyclopentadiene			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4,6-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4,5-trichlorophenol			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-chloronaphthalene			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
acenaphthylene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
dimethylphthalate			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,6-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4-dinitrotoluene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
acenaphthene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
3-nitroaniline			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4-dinitrophenol			< 5	5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
dibenzofuran			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-nitrophenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
fluorene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
diethyl phthalate			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12 9:30	Reporting Result	Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis Date	Time	Reference
4-chlorophenyl phenyl ether			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-nitroaniline			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4,6-dinitro-2-methylphenol			< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
azobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
N-nitrosodiphenylamine			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
4-bromophenyl phenyl ether			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
hexachlorobenzene			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
pentachlorophenol			< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
phenanthrene			0.21	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
anthracene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
carbazole			< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
di-n-butylphthalate			< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
fluoranthene			0.21	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzidine			< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
pyrene			0.18	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
butyl benzyl phthalate			0.6	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzo(a)anthracene			0.06	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
chrysene			0.09	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
3,3'-dichlorobenzidine			< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
bis(2-ethylhexyl)phthalate			2.4	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
di-n-octyl phthalate			0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzo(b)fluoranthene			0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzo(k)fluoranthene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzo(a)pyrene			0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
indeno(1,2,3-cd)pyrene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
dibenzo(a,h)anthracene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
benzo(g,h,i)perylene			< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
Surrogate Recovery												
Limits												
2-fluorophenol SUR			2 *	21-100	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
phenol-D5 SUR			18	10-102	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2,4,6-tribromophenol SUR			0 *	10-123	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
nitrobenzene-D5 SUR			42	35-114	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
2-fluorobiphenyl SUR			66	43-116	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D
p-terphenyl-D14 SUR			55	33-141	%	1	AJD	10/10/12	5556	10/12/12	21:08	SW3546/8270D

* This surrogate showed recovery outside the acceptance limits as a result of interferences caused by large co-eluting non-target compounds.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Sampled:	10/3/12 9:58	Reporting	Instr	Dil'n	Prep	Analysis					
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
N-nitrosodimethylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
aniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
phenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-chlorophenol		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
bis(2-chloroethyl)ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
1,3-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
1,4-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
1,2-dichlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzyl alcohol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
bis(2-chloroisopropyl) ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
hexachloroethane		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
N-nitroso-di-N-propylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
nitrobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
isophorone		1.8	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-nitrophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4-dimethylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
bis(2-chloroethoxy)methane		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4-dichlorophenol		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
1,2,4-trichlorobenzene		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
naphthalene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzoic acid		< 5	5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-chloroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
hexachlorobutadiene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-chloro-3-methylphenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-methylnaphthalene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
hexachlorocyclopentadiene		< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4,6-trichlorophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4,5-trichlorophenol		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-chloronaphthalene		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-nitroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
acenaphthylene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
dimethylphthalate		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,6-dinitrotoluene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4-dinitrotoluene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
acenaphthene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
3-nitroaniline		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4-dinitrophenol		< 5	5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
dibenzofuran		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-nitrophenol		< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
fluorene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
diethyl phthalate		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:58	Reporting	Instr	Dil'n	Prep	Analysis					
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether	< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-nitroaniline	< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4,6-dinitro-2-methylphenol	< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
azobenzene	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
N-nitrosodiphenylamine	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
4-bromophenyl phenyl ether	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
hexachlorobenzene	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
pentachlorophenol	< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
phenanthrene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
anthracene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
carbazole	< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
di-n-butylphthalate	< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
fluoranthene	0.07	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzidine	< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
pyrene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
butyl benzyl phthalate	< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzo(a)anthracene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
chrysene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
3,3'-dichlorobenzidine	< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
bis(2-ethylhexyl)phthalate	0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
di-n-octyl phthalate	< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzo(b)fluoranthene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzo(k)fluoranthene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzo(a)pyrene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
indeno(1,2,3-cd)pyrene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
dibenzo(a,h)anthracene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
benzo(g,h,i)perylene	< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
Surrogate Recovery										
Limits										
2-fluorophenol SUR	0 *	21-100	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
phenol-D5 SUR	3 *	10-102	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2,4,6-tribromophenol SUR	0 *	10-123	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
nitrobenzene-D5 SUR	39	35-114	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
2-fluorobiphenyl SUR	52	43-116	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D
p-terphenyl-D14 SUR	53	33-141	%	1	AJD	10/10/12	5556	10/12/12	14:18	SW3546/8270D

* This surrogate showed recovery outside the acceptance limits as a result of interferences caused by large co-eluting non-target compounds.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Parameter	Sampled:	10/3/12	10:40	Reporting	Instr	Dil'n	Prep	Analysis			Reference	
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	
N-nitrosodimethylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
aniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
phenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-chlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
bis(2-chloroethyl)ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
1,3-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
1,4-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
1,2-dichlorobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzyl alcohol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
bis(2-chloroisopropyl) ether				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
hexachloroethane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
N-nitroso-di-N-propylamine				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
nitrobenzene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
isophorone				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-nitrophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4-dimethylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
bis(2-chloroethoxy)methane				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4-dichlorophenol				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
1,2,4-trichlorobenzene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
naphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzoic acid				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-chloroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
hexachlorobutadiene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-chloro-3-methylphenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-methylnaphthalene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
hexachlorocyclopentadiene				< 1	1	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4,6-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4,5-trichlorophenol				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-chloronaphthalene				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
acenaphthylene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
dimethylphthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,6-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4-dinitrotoluene				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
acenaphthene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
3-nitroaniline				< 0.2	0.2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4-dinitrophenol				< 5	5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
dibenzofuran				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-nitrophenol				< 2	2	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
fluorene				< 0.05	0.05	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D
diethyl phthalate				< 0.5	0.5	ug/g	1	AJD 10/10/12	5556	10/12/12	14:56	SW3546/8270D

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Sampled:	10/3/12 10:40	Reporting	Instr	Dil'n	Prep	Analysis					
Parameter		Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
4-chlorophenyl phenyl ether		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-nitroaniline		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
4,6-dinitro-2-methylphenol		< 2	2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
azobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
N-nitrosodiphenylamine		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
4-bromophenyl phenyl ether		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
hexachlorobenzene		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
pentachlorophenol		< 1	1	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
phenanthrene		0.19	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
anthracene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
carbazole		< 0.2	0.2	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
di-n-butylphthalate		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
fluoranthene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzidine		< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
pyrene		0.09	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
butyl benzyl phthalate		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzo(a)anthracene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
chrysene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
3,3'-dichlorobenzidine		< 3	3	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
bis(2-ethylhexyl)phthalate		2.6	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
di-n-octyl phthalate		< 0.5	0.5	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzo(b)fluoranthene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzo(k)fluoranthene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzo(a)pyrene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
indeno(1,2,3-cd)pyrene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
dibenzo(a,h)anthracene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
benzo(g,h,i)perylene		< 0.05	0.05	ug/g	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
Surrogate Recovery											
Limits											
2-fluorophenol SUR		4 *	21-100	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
phenol-D5 SUR		15	10-102	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
2,4,6-tribromophenol SUR		0 *	10-123	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
nitrobenzene-D5 SUR		51	35-114	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
2-fluorobiphenyl SUR		71	43-116	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D
p-terphenyl-D14 SUR		61	33-141	%	1	AJD	10/10/12	5556	10/12/12	14:56	SW3546/8270D

* This surrogate showed recovery outside the acceptance limits as a result of interferences caused by large co-eluting non-target compounds.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-003

Sample ID: CA-SB-3 10-12.5

Matrix: Solid

Percent Dry: 72.3% Results expressed on a dry weight basis.

Sampled: 10/2/12 13:10		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	60	30-150	%	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A
decachlorobiphenyl SUR	71	30-150	%	1	JLZ	10/8/12	5551	10/9/12	19:13	SW3540C8082A

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Sampled: 10/3/12 13:30		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1221	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1232	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1242	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1248	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1254	< 7.1	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
PCB-1260	250	7.1	ug/g	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	DOR	30-150	%	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A
decachlorobiphenyl SUR	DOR	30-150	%	200	JLZ	10/8/12	5551	10/12/12	5:05	SW3540C8082A

DOR = Diluted out of range.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-006

Sample ID: CA-SB-6 0-5

Matrix: Solid

Percent Dry: 93.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	71	30-150	%	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A
decachlorobiphenyl SUR	74	30-150	%	1	JLZ	10/8/12	5551	10/11/12	18:57	SW3540C8082A

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:50		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	71	30-150	%	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A
decachlorobiphenyl SUR	82	30-150	%	1	JLZ	10/8/12	5551	10/11/12	19:27	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Sampled: 10/2/12 16:15		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	69	30-150	%	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A
decachlorobiphenyl SUR	75	30-150	%	1	JLZ	10/8/12	5551	10/9/12	21:14	SW3540C8082A

Sample#: 25123-009

Sample ID: CA-SB-8 0-2.5

Matrix: Solid

Percent Dry: 97.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:50		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	65	30-150	%	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A
decachlorobiphenyl SUR	80	30-150	%	1	JLZ	10/8/12	5551	10/9/12	21:45	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-010

Sample ID: CA-SB-8 5-10

Matrix: Solid

Percent Dry: 83.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 10:00		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	62	30-150	%	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A
decachlorobiphenyl SUR	61	30-150	%	1	JLZ	10/9/12	5551	10/11/12	19:58	SW3540C8082A

Sample#: 25123-011

Sample ID: CA-SB-9 0-4

Matrix: Solid

Percent Dry: 82.4% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:10		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	68	30-150	%	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A
decachlorobiphenyl SUR	74	30-150	%	1	JLZ	10/9/12	5551	10/11/12	20:28	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:30		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	59	30-150	%	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	
decachlorobiphenyl SUR	58	30-150	%	1	JLZ	10/10/12	5561	10/12/12	1:32	SW3540C8082A	

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:30		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	72	30-150	%	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	
decachlorobiphenyl SUR	74	30-150	%	1	JLZ	10/9/12	5551	10/12/12	13:43	SW3540C8082A	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-015

Sample ID: CA-SB-12 15-18

Matrix: Solid

Percent Dry: 52.1% Results expressed on a dry weight basis.

Sampled: 10/2/12 8:45		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1221	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1232	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1242	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1248	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1254	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
PCB-1260	< 0.06	0.06	ug/g	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	62	30-150	%	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A
decachlorobiphenyl SUR	75	30-150	%	1	JLZ	10/9/12	5551	10/11/12	21:29	SW3540C8082A

Sample#: 25123-016

Sample ID: CA-SB-13 10-12

Matrix: Solid

Percent Dry: 84.2% Results expressed on a dry weight basis.

Sampled: 10/2/12 9:10		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	54	30-150	%	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A
decachlorobiphenyl SUR	52	30-150	%	1	JLZ	10/9/12	5551	10/11/12	22:00	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-017

Sample ID: SB-DUP-1

Matrix: Solid

Percent Dry: 87.8% Results expressed on a dry weight basis.

Sampled: 10/2/12 0:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	51	30-150	%	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A
decachlorobiphenyl SUR	58	30-150	%	1	JLZ	10/9/12	5551	10/11/12	22:30	SW3540C8082A

Sample#: 25123-020

Sample ID: CA-SS-1

Matrix: Solid

Percent Dry: 82.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:20		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	66	30-150	%	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A
decachlorobiphenyl SUR	75	30-150	%	1	JLZ	10/9/12	5551	10/11/12	23:00	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-021

Sample ID: CA-SS-2

Matrix: Solid

Percent Dry: 91.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:10		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	60	30-150	%	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A
decachlorobiphenyl SUR	62	30-150	%	1	JLZ	10/9/12	5551	10/11/12	23:31	SW3540C8082A

Sample#: 25123-022

Sample ID: CA-SS-3

Matrix: Solid

Percent Dry: 93.4% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:15		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	69	30-150	%	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A
decachlorobiphenyl SUR	75	30-150	%	1	JLZ	10/9/12	5551	10/12/12	1:02	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-023

Sample ID: CA-SS-4

Matrix: Solid

Percent Dry: 91.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:45		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
PCB-1260	0.75	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	65	30-150	%	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	
decachlorobiphenyl SUR	72	30-150	%	1	JLZ	10/10/12	5561	10/12/12	2:03	SW3540C8082A	

Sample#: 25123-024

Sample ID: CA-SS-5

Matrix: Solid

Percent Dry: 81.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 13:00		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	61	30-150	%	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	
decachlorobiphenyl SUR	57	30-150	%	1	JLZ	10/10/12	5561	10/12/12	2:33	SW3540C8082A	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:30		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1221	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1232	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1242	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1248	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1254	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
PCB-1260	< 0.1	0.1	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	82	30-150	%	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A
decachlorobiphenyl SUR	58	30-150	%	1	JLZ	10/10/12	5561	10/12/12	3:03	SW3540C8082A

Note: Elevated reporting limit due to matrix interference.

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:58		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1221	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1232	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1242	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1248	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1254	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
PCB-1260	< 0.2	0.2	ug/g	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	90	30-150	%	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A
decachlorobiphenyl SUR	97	30-150	%	1	JLZ	10/10/12	5561	10/12/12	3:34	SW3540C8082A

Note: Elevated reporting limit due to matrix interference.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Sampled: 10/3/12 10:40		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	57	30-150	%	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A
decachlorobiphenyl SUR	62	30-150	%	1	JLZ	10/10/12	5561	10/12/12	4:04		SW3540C8082A

Sample#: 25123-028

Sample ID: CA-CC-4

Matrix: Solid

Percent Dry: 96.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:25		Reporting		Instr	Dil'n	Prep		Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
Surrogate Recovery											
Limits											
tetrachloro-m-xylene SUR	40	30-150	%	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A
decachlorobiphenyl SUR	40	30-150	%	1	JLZ	10/10/12	5561	10/12/12	4:34		SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-029

Sample ID: CA-CC-5

Matrix: Solid

Percent Dry: 96.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:48

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1221	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1232	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1242	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1248	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1254	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
PCB-1260	< 0.2	0.2	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	#	30-150	%	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A
decachlorobiphenyl SUR	139	30-150	%	1	JLZ	10/11/12	5561	10/12/12	10:13	SW3540C8082A

Note: Elevated reporting limit due to matrix interference.

This surrogate is above the acceptance criteria. Since no targets were detected above the quantitation limit, there is no impact to the data.

Sample#: 25123-030

Sample ID: STACK

Matrix: Solid

Percent Dry: 99.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 14:50

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	61	30-150	%	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A
decachlorobiphenyl SUR	77	30-150	%	1	JLZ	10/11/12	5561	10/12/12	10:44	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-002

Sample ID: CA-SB-2 0-1

Matrix: Solid

Percent Dry: 87.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	15:00	Reporting	Instr	Dil'n	Prep	Analysis					
				Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene				0.7	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
2-methylnaphthalene				0.8	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
phenanthrene				22	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
acenaphthene				3.0	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
acenaphthylene				< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
fluorene				3.3	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
anthracene				6.0	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
fluoranthene				15	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
pyrene				15	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
benzo(a)anthracene				8.1	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
chrysene				7.3	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
benzo(b)fluoranthene				4.2	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
benzo(k)fluoranthene				5.2	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
benzo(a)pyrene				5.3	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
indeno(1,2,3-cd)pyrene				1.7	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
dibenzo(a,h)anthracene				1.1	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
benzo(g,h,i)perylene				1.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	22:39	MA EPH
Unadjusted C11-C22 Aromatics				180	23	ug/g	1	JLZ	10/8/12	5547	10/9/12	1:18	MA EPH
C9-C18 Aliphatics				< 23	23	ug/g	1	JLZ	10/8/12	5547	10/9/12	20:27	MA EPH
C19-C36 Aliphatics				< 23	23	ug/g	1	JLZ	10/8/12	5547	10/9/12	20:27	MA EPH
C11-C22 Aromatics				79	23	ug/g	1	JLZ	10/8/12	5547	10/9/12	1:18	MA EPH
Surrogate Recovery													
Limits													
1-chloro-octadecane SUR				38 *	40-140	%	1	JLZ	10/8/12	5547	10/9/12	20:27	MA EPH
o-terphenyl SUR				31 *	40-140	%	1	JLZ	10/8/12	5547	10/9/12	1:18	MA EPH
2-fluorobiphenyl SUR				69	40-140	%	1	JLZ	10/8/12	5547	10/9/12	1:18	MA EPH
2-bromonaphthalene SUR				74	40-140	%	1	JLZ	10/8/12	5547	10/9/12	1:18	MA EPH

* The surrogate showed recovery outside the acceptance limits. Results from re-extraction did not compare well. Both sets of data are provided.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-003

Sample ID: CA-SB-3 10-12.5

Matrix: Solid

Percent Dry: 72.3% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	13:10	Reporting	Instr	Dil'n	Prep	Analysis				
				Result	Limit	Units	Analyst	Date	Date	Time	Reference	
naphthalene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
2-methylnaphthalene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
phenanthrene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
acenaphthene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
acenaphthylene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
fluorene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
anthracene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
fluoranthene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
pyrene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
benzo(a)anthracene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
chrysene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
benzo(b)fluoranthene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
benzo(k)fluoranthene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
benzo(a)pyrene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
indeno(1,2,3-cd)pyrene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
dibenzo(a,h)anthracene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
benzo(g,h,i)perylene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	14:22	MA EPH
Unadjusted C11-C22 Aromatics				< 27	27	ug/g	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
C9-C18 Aliphatics				< 27	27	ug/g	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
C19-C36 Aliphatics				< 27	27	ug/g	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
C11-C22 Aromatics				< 27	27	ug/g	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
Surrogate Recovery												
						Limits						
1-chloro-octadecane SUR				50	40-140	%	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
o-terphenyl SUR				69	40-140	%	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
2-fluorobiphenyl SUR				N/A	40-140	%	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH
2-bromonaphthalene SUR				N/A	40-140	%	1	JLZ 10/8/12	5547	10/15/12	15:30	MA EPH

Note: No ranges were detected above the method reporting limit; therefore fractionation was not required.

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	11:50	Reporting	Instr	Dil'n	Prep	Analysis				
				Result	Limit	Units	Analyst	Date	Date	Time	Reference	
Unadjusted C11-C22 Aromatics				< 20	20	ug/g	1	JLZ 10/16/12	5572	10/17/12	15:16	MA EPH
C9-C18 Aliphatics				< 20	20	ug/g	1	JLZ 10/16/12	5572	10/16/12	20:34	MA EPH
C19-C36 Aliphatics				< 20	20	ug/g	1	JLZ 10/16/12	5572	10/16/12	20:34	MA EPH
Surrogate Recovery												
						Limits						
1-chloro-octadecane SUR				62	40-140	%	1	JLZ 10/16/12	5572	10/16/12	20:34	MA EPH
o-terphenyl SUR				70	40-140	%	1	JLZ 10/16/12	5572	10/17/12	15:16	MA EPH
2-fluorobiphenyl SUR				105	40-140	%	1	JLZ 10/16/12	5572	10/17/12	15:16	MA EPH
2-bromonaphthalene SUR				108	40-140	%	1	JLZ 10/16/12	5572	10/17/12	15:16	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:50		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Unadjusted C11-C22 Aromatics	< 24	24	ug/g	1	JLZ	10/8/12	5547	10/14/12	20:12	MA EPH
C9-C18 Aliphatics	< 24	24	ug/g	1	JLZ	10/8/12	5547	10/9/12	21:21	MA EPH
C19-C36 Aliphatics	69	24	ug/g	1	JLZ	10/8/12	5547	10/9/12	21:21	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	52	40-140	%	1	JLZ	10/8/12	5547	10/9/12	21:21	MA EPH
o-terphenyl SUR	46	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:12	MA EPH
2-fluorobiphenyl SUR	76	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:12	MA EPH
2-bromonaphthalene SUR	82	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:12	MA EPH

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Sampled: 10/2/12 16:15		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Unadjusted C11-C22 Aromatics	< 22	22	ug/g	1	JLZ	10/8/12	5547	10/14/12	20:35	MA EPH
C9-C18 Aliphatics	< 22	22	ug/g	1	JLZ	10/8/12	5547	10/9/12	21:40	MA EPH
C19-C36 Aliphatics	< 22	22	ug/g	1	JLZ	10/8/12	5547	10/9/12	21:40	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	50	40-140	%	1	JLZ	10/8/12	5547	10/9/12	21:40	MA EPH
o-terphenyl SUR	45	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:35	MA EPH
2-fluorobiphenyl SUR	74	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:35	MA EPH
2-bromonaphthalene SUR	77	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:35	MA EPH

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:30		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
Unadjusted C11-C22 Aromatics	< 25	25	ug/g	1	JLZ	10/8/12	5547	10/14/12	20:57	MA EPH
C9-C18 Aliphatics	< 25	25	ug/g	1	JLZ	10/8/12	5547	10/9/12	23:29	MA EPH
C19-C36 Aliphatics	88	25	ug/g	1	JLZ	10/8/12	5547	10/9/12	23:29	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	48	40-140	%	1	JLZ	10/8/12	5547	10/9/12	23:29	MA EPH
o-terphenyl SUR	50	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:57	MA EPH
2-fluorobiphenyl SUR	87	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:57	MA EPH
2-bromonaphthalene SUR	94	40-140	%	1	JLZ	10/8/12	5547	10/14/12	20:57	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-014

Sample ID: SB-DUP-2

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 10/3/12 0:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Unadjusted C11-C22 Aromatics	< 25	25	ug/g	1	JLZ	10/8/12	5547	10/14/12	21:20	MA EPH
C9-C18 Aliphatics	< 25	25	ug/g	1	JLZ	10/8/12	5547	10/9/12	23:47	MA EPH
C19-C36 Aliphatics	36	25	ug/g	1	JLZ	10/8/12	5547	10/9/12	23:47	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	51	40-140	%	1	JLZ	10/8/12	5547	10/9/12	23:47	MA EPH
o-terphenyl SUR	42	40-140	%	1	JLZ	10/8/12	5547	10/14/12	21:20	MA EPH
2-fluorobiphenyl SUR	72	40-140	%	1	JLZ	10/8/12	5547	10/14/12	21:20	MA EPH
2-bromonaphthalene SUR	71	40-140	%	1	JLZ	10/8/12	5547	10/14/12	21:20	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-015

Sample ID: CA-SB-12 15-18

Matrix: Solid

Percent Dry: 52.1% Results expressed on a dry weight basis.

Sampled: 10/2/12 8:45		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene	0.4	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
2-methylnaphthalene	1.1	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
phenanthrene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
acenaphthene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
acenaphthylene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
fluorene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
anthracene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
fluoranthene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
pyrene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
benzo(a)anthracene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
chrysene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
benzo(b)fluoranthene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
benzo(k)fluoranthene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
benzo(a)pyrene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
indeno(1,2,3-cd)pyrene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
dibenzo(a,h)anthracene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
benzo(g,h,i)perylene	< 0.2	0.2	ug/g	1	AJD	10/8/12	5547	10/9/12	13:45	MA EPH
Unadjusted C11-C22 Aromatics	340	35	ug/g	1	JLZ	10/8/12	5547	10/14/12	22:28	MA EPH
C9-C18 Aliphatics	990	35	ug/g	1	JLZ	10/8/12	5547	10/10/12	0:06	MA EPH
C19-C36 Aliphatics	310	35	ug/g	1	JLZ	10/8/12	5547	10/10/12	0:06	MA EPH
C11-C22 Aromatics	340	35	ug/g	1	JLZ	10/8/12	5547	10/14/12	22:28	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	19 *	40-140	%	1	JLZ	10/8/12	5547	10/10/12	0:06	MA EPH
o-terphenyl SUR	33 *	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:28	MA EPH
2-fluorobiphenyl SUR	101	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:28	MA EPH
2-bromonaphthalene SUR	64	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:28	MA EPH

* The surrogate showed recovery outside the acceptance limits as a result of hydrocarbons present in the sample. The chromatogram is included for reference.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-016

Sample ID: CA-SB-13 10-12

Matrix: Solid

Percent Dry: 84.2% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12 9:10	Reporting	Instr	Dil'n	Prep	Analysis					
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
2-methylnaphthalene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
phenanthrene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
acenaphthene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
acenaphthylene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
fluorene			0.7	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
anthracene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
fluoranthene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
pyrene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
benzo(a)anthracene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
chrysene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
benzo(b)fluoranthene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
benzo(k)fluoranthene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
benzo(a)pyrene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
indeno(1,2,3-cd)pyrene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
dibenzo(a,h)anthracene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
benzo(g,h,i)perylene			< 0.6	0.6	ug/g	5	AJD	10/8/12	5547	10/11/12	21:25	MA EPH
Unadjusted C11-C22 Aromatics			1300	23	ug/g	1	JLZ	10/8/12	5547	10/14/12	22:51	MA EPH
C9-C18 Aliphatics			1600	23	ug/g	1	JLZ	10/8/12	5547	10/10/12	1:00	MA EPH
C19-C36 Aliphatics			1900	23	ug/g	1	JLZ	10/8/12	5547	10/10/12	1:00	MA EPH
C11-C22 Aromatics			1300	23	ug/g	1	JLZ	10/8/12	5547	10/14/12	22:51	MA EPH
Surrogate Recovery												
Limits												
1-chloro-octadecane SUR			#	40-140	%	1	JLZ	10/8/12	5547	10/10/12	1:00	MA EPH
o-terphenyl SUR			41	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:51	MA EPH
2-fluorobiphenyl SUR			109	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:51	MA EPH
2-bromonaphthalene SUR			72	40-140	%	1	JLZ	10/8/12	5547	10/14/12	22:51	MA EPH

The surrogate could not be distinguished from hydrocarbons present in the chromatogram. The chromatogram is included for reference.

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-017

Sample ID: SB-DUP-1

Matrix: Solid

Percent Dry: 87.8% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12 0:00	Reporting	Instr	Dil'n	Prep	Analysis					
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
2-methylnaphthalene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
phenanthrene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
acenaphthene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
acenaphthylene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
fluorene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
anthracene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
fluoranthene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
pyrene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
benzo(a)anthracene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
chrysene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
benzo(b)fluoranthene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
benzo(k)fluoranthene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
benzo(a)pyrene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
indeno(1,2,3-cd)pyrene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
dibenzo(a,h)anthracene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
benzo(g,h,i)perylene			< 0.5	0.5	ug/g	5	AJD	10/8/12	5547	10/11/12	22:02	MA EPH
Unadjusted C11-C22 Aromatics			670	22	ug/g	1	JLZ	10/8/12	5547	10/14/12	23:13	MA EPH
C9-C18 Aliphatics			1100	22	ug/g	1	JLZ	10/8/12	5547	10/10/12	1:19	MA EPH
C19-C36 Aliphatics			1400	22	ug/g	1	JLZ	10/8/12	5547	10/10/12	1:19	MA EPH
C11-C22 Aromatics			670	22	ug/g	1	JLZ	10/8/12	5547	10/14/12	23:13	MA EPH
Surrogate Recovery												
Limits												
1-chloro-octadecane SUR			#	40-140	%	1	JLZ	10/8/12	5547	10/10/12	1:19	MA EPH
o-terphenyl SUR			35 *	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:13	MA EPH
2-fluorobiphenyl SUR			91	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:13	MA EPH
2-bromonaphthalene SUR			65	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:13	MA EPH

The surrogate could not be distinguished from hydrocarbons present in the chromatogram. The chromatogram is included for reference.

*** The surrogate showed recovery outside the acceptance limits as a result of hydrocarbons present in the sample.**

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-018

Sample ID: CA-SB-14 0-3

Matrix: Solid

Percent Dry: 94.8% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12	10:15	Reporting	Instr	Dil'n	Prep	Analysis				
				Result	Limit	Units	Analyst	Date	Batch	Date	Time	Reference
naphthalene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
2-methylnaphthalene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
phenanthrene				0.4	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
acenaphthene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
acenaphthylene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
fluorene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
anthracene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
fluoranthene				1.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
pyrene				1.0	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
benzo(a)anthracene				0.7	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
chrysene				0.6	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
benzo(b)fluoranthene				0.6	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
benzo(k)fluoranthene				0.8	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
benzo(a)pyrene				0.6	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
indeno(1,2,3-cd)pyrene				0.2	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
dibenzo(a,h)anthracene				< 0.1	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
benzo(g,h,i)perylene				0.2	0.1	ug/g	1	AJD 10/8/12	5547	10/9/12	17:29	MA EPH
Unadjusted C11-C22 Aromatics				34	20	ug/g	1	JLZ 10/8/12	5547	10/14/12	23:36	MA EPH
C9-C18 Aliphatics				< 20	20	ug/g	1	JLZ 10/8/12	5547	10/10/12	1:37	MA EPH
C19-C36 Aliphatics				58	20	ug/g	1	JLZ 10/8/12	5547	10/10/12	1:37	MA EPH
C11-C22 Aromatics				27	20	ug/g	1	JLZ 10/8/12	5547	10/14/12	23:36	MA EPH
Surrogate Recovery												
Limits												
1-chloro-octadecane SUR				46	40-140	%	1	JLZ 10/8/12	5547	10/10/12	1:37	MA EPH
o-terphenyl SUR				48	40-140	%	1	JLZ 10/8/12	5547	10/14/12	23:36	MA EPH
2-fluorobiphenyl SUR				77	40-140	%	1	JLZ 10/8/12	5547	10/14/12	23:36	MA EPH
2-bromonaphthalene SUR				83	40-140	%	1	JLZ 10/8/12	5547	10/14/12	23:36	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-019

Sample ID: CA-SB-15 5-5.5

Matrix: Solid

Percent Dry: 88.1% Results expressed on a dry weight basis.

Sampled: 10/2/12 10:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
naphthalene	< 0.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
2-methylnaphthalene	< 0.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
phenanthrene	1.5	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
acenaphthene	0.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
acenaphthylene	< 0.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
fluorene	0.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
anthracene	0.3	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
fluoranthene	2.7	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
pyrene	2.2	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
benzo(a)anthracene	1.2	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
chrysene	1.3	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
benzo(b)fluoranthene	1.1	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
benzo(k)fluoranthene	1.3	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
benzo(a)pyrene	1.2	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
indeno(1,2,3-cd)pyrene	0.4	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
dibenzo(a,h)anthracene	0.2	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
benzo(g,h,i)perylene	0.4	0.1	ug/g	1	AJD	10/8/12	5547	10/9/12	15:37	MA EPH
Unadjusted C11-C22 Aromatics	32	21	ug/g	1	JLZ	10/8/12	5547	10/14/12	23:58	MA EPH
C9-C18 Aliphatics	< 21	21	ug/g	1	JLZ	10/8/12	5547	10/10/12	0:24	MA EPH
C19-C36 Aliphatics	< 21	21	ug/g	1	JLZ	10/8/12	5547	10/10/12	0:24	MA EPH
C11-C22 Aromatics	< 21	21	ug/g	1	JLZ	10/8/12	5547	10/14/12	23:58	MA EPH
Surrogate Recovery										
Limits										
1-chloro-octadecane SUR	55	40-140	%	1	JLZ	10/8/12	5547	10/10/12	0:24	MA EPH
o-terphenyl SUR	51	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:58	MA EPH
2-fluorobiphenyl SUR	77	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:58	MA EPH
2-bromonaphthalene SUR	84	40-140	%	1	JLZ	10/8/12	5547	10/14/12	23:58	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-032

Sample ID: CA-SB-2 0-1 Re-Extraction

Matrix: Solid

Percent Dry: 87.6% Results expressed on a dry weight basis.

Parameter	Sampled:	10/2/12 15:00	Reporting	Instr	Dil'n	Prep	Analysis					
			Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
naphthalene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
2-methylnaphthalene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
phenanthrene			0.8	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
acenaphthene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
acenaphthylene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
fluorene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
anthracene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
fluoranthene			0.7	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
pyrene			0.9	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
benzo(a)anthracene			0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
chrysene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
benzo(b)fluoranthene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
benzo(k)fluoranthene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
benzo(a)pyrene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
indeno(1,2,3-cd)pyrene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
dibenzo(a,h)anthracene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
benzo(g,h,i)perylene			< 0.6	0.6	ug/g	5	AJD	10/16/12	5572	10/17/12	11:42	MA EPH
Unadjusted C11-C22 Aromatics			< 22	22	ug/g	1	JLZ	10/16/12	5572	10/17/12	14:53	MA EPH
C9-C18 Aliphatics			< 22	22	ug/g	1	JLZ	10/16/12	5572	10/16/12	20:16	MA EPH
C19-C36 Aliphatics			< 22	22	ug/g	1	JLZ	10/16/12	5572	10/16/12	20:16	MA EPH
C11-C22 Aromatics			< 22	22	ug/g	1	JLZ	10/16/12	5572	10/17/12	14:53	MA EPH
Surrogate Recovery												
Limits												
1-chloro-octadecane SUR			45	40-140	%	1	JLZ	10/16/12	5572	10/16/12	20:16	MA EPH
o-terphenyl SUR			60	40-140	%	1	JLZ	10/16/12	5572	10/17/12	14:53	MA EPH
2-fluorobiphenyl SUR			112	40-140	%	1	JLZ	10/16/12	5572	10/17/12	14:53	MA EPH
2-bromonaphthalene SUR			118	40-140	%	1	JLZ	10/16/12	5572	10/17/12	14:53	MA EPH

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-001

Sample ID: CA-SB-1 2.5-5

Matrix: Solid Percent Dry: 84.1% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 15:30		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Lead	100	0.7	ug/g	1	SAS	10/8/12	5553	10/9/12	23:56	SW3051A6010C	

Sample#: 25123-002

Sample ID: CA-SB-2 0-1

Matrix: Solid Percent Dry: 87.6% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 15:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Antimony	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Arsenic	14	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Beryllium	0.8	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Chromium	25	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Copper	37	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Lead	210	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Mercury	0.44	0.16	ug/g	1	SAS	10/12/12	5564	10/12/12	16:29	SW7471B	
Nickel	18	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Thallium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	
Zinc	80	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:07	SW3051A6010C	

Sample#: 25123-003

Sample ID: CA-SB-3 10-12.5

Matrix: Solid Percent Dry: 72.3% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 13:10		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	11	0.8	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Barium	85	4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Chromium	51	4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Lead	11	0.8	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Mercury	< 0.20	0.20	ug/g	1	SAS	10/12/12	5564	10/12/12	16:31	SW7471B	
Selenium	< 4	4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	
Silver	< 0.5	0.5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:03	SW3051A6010C	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-004

Sample ID: CA-SB-4 0-2

Matrix: Solid

Percent Dry: 97.2% Results expressed on a dry weight basis.

Sampled: 10/2/12 11:50

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Arsenic	4.8	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Beryllium	0.3	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Chromium	13	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Copper	10	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Lead	5.8	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5564	10/12/12	16:33	SW7471B
Nickel	11	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C
Zinc	20	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:21	SW3051A6010C

Sample#: 25123-005

Sample ID: CA-SB-5 5-10

Matrix: Solid

Percent Dry: 85.6% Results expressed on a dry weight basis.

Sampled: 10/3/12 13:30

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Arsenic	17	0.5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Barium	70	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Chromium	22	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Lead	79	0.5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Mercury	< 0.17	0.17	ug/g	1	SAS	10/12/12	5564	10/12/12	16:35	SW7471B
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:10	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-006

Sample ID: CA-SB-6 0-5

Matrix: Solid

Percent Dry: 93.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Arsenic	7.1	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Beryllium	0.4	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Chromium	13	2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Copper	15	2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Lead	23	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5564	10/12/12	16:37	SW7471B
Nickel	11	2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Selenium	< 2	2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Silver	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C
Zinc	36	2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:29	SW3051A6010C

Sample#: 25123-007

Sample ID: CA-SB-6 12.5-15

Matrix: Solid

Percent Dry: 77% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:50		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
Antimony	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Arsenic	8.7	0.7	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Beryllium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Chromium	19	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Copper	43	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Lead	22	0.7	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Mercury	< 0.19	0.19	ug/g	1	SAS	10/12/12	5564	10/12/12	16:38	SW7471B
Nickel	13	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Silver	< 0.5	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Thallium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C
Zinc	32	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:36	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-008

Sample ID: CA-SB-7 7.5-10

Matrix: Solid

Percent Dry: 88.4% Results expressed on a dry weight basis.

Sampled: 10/2/12 16:15

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Arsenic	3.2	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Beryllium	0.5	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Chromium	12	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Copper	25	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Lead	15	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Mercury	< 0.16	0.16	ug/g	1	SAS	10/12/12	5564	10/12/12	16:40	SW7471B
Nickel	6	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C
Zinc	30	3	ug/g	1	SAS	10/8/12	5553	10/9/12	21:44	SW3051A6010C

Sample#: 25123-009

Sample ID: CA-SB-8 0-2.5

Matrix: Solid

Percent Dry: 97.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 8:50

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Arsenic	11	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Beryllium	0.3	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Chromium	7	2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Copper	11	2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Lead	3.4	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5564	10/12/12	16:42	SW7471B
Nickel	7	2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Selenium	< 2	2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Silver	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C
Zinc	13	2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:20	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-010

Sample ID: CA-SB-8 5-10

Matrix: Solid

Percent Dry: 83.8% Results expressed on a dry weight basis.

Parameter	Sampled: 10/3/12 10:00		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Antimony	16	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Arsenic	18	0.7	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Beryllium	0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Chromium	30	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Copper	990	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Lead	540	0.7	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Mercury	0.27	0.17	ug/g	1	SAS	10/12/12	5564	10/12/12	16:55	SW7471B	
Nickel	44	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Silver	0.5	0.5	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Thallium	< 0.3	0.3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	
Zinc	360	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:27	SW3051A6010C	

Sample#: 25123-012

Sample ID: CA-SB-10 10-15

Matrix: Solid

Percent Dry: 79.7% Results expressed on a dry weight basis.

Parameter	Sampled: 10/3/12 12:30		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	11	0.6	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Barium	59	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Chromium	27	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Lead	29	0.6	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Mercury	< 0.18	0.18	ug/g	1	SAS	10/12/12	5564	10/12/12	16:58	SW7471B	
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:26	SW3051A6010C	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-013

Sample ID: CA-SB-11 10-15

Matrix: Solid

Percent Dry: 78.5% Results expressed on a dry weight basis.

Parameter	Sampled: 10/3/12 8:30		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Antimony	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Arsenic	9.7	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Beryllium	1.0	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Cadmium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Chromium	31	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Copper	50	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Lead	75	0.6	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Mercury	0.43	0.18	ug/g	1	SAS	10/12/12	5564	10/12/12	17:00	SW7471B	
Nickel	24	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Selenium	< 3	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Silver	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Thallium	< 0.2	0.2	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	
Zinc	68	3	ug/g	1	SAS	10/8/12	5553	10/9/12	22:35	SW3051A6010C	

Sample#: 25123-015

Sample ID: CA-SB-12 15-18

Matrix: Solid

Percent Dry: 52.1% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 8:45		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	8.4	1.0	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Barium	43	5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Cadmium	< 0.4	0.4	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Chromium	15	5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Lead	320	1.0	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Mercury	0.39	0.28	ug/g	1	SAS	10/12/12	5564	10/12/12	17:02	SW7471B	
Selenium	< 5	5	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	
Silver	< 0.7	0.7	ug/g	1	SAS	10/8/12	5553	10/10/12	0:41	SW3051A6010C	

Sample#: 25123-016

Sample ID: CA-SB-13 10-12

Matrix: Solid

Percent Dry: 84.2% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 9:10		Reporting		Instr	Dil'n	Prep	Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	3.2	0.7	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Barium	22	3	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Chromium	11	3	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Lead	5.2	0.7	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Mercury	< 0.17	0.17	ug/g	1	SAS	10/12/12	5564	10/12/12	17:04	SW7471B	
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	
Silver	< 0.5	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	22:57	SW3051A6010C	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-017

Sample ID: SB-DUP-1

Matrix: Solid

Percent Dry: 87.8% Results expressed on a dry weight basis.

Parameter	Sampled: 10/2/12 0:00		Reporting		Instr Dil'n	Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Arsenic	3.4	0.7	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Barium	28	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Cadmium	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Chromium	12	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Lead	7.1	0.7	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Mercury	< 0.16	0.16	ug/g	1	SAS	10/12/12	5564	10/12/12	17:06	SW7471B	
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	
Silver	< 0.5	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	23:11	SW3051A6010C	

Sample#: 25123-020

Sample ID: CA-SS-1

Matrix: Solid

Percent Dry: 82.2% Results expressed on a dry weight basis.

Parameter	Sampled: 10/3/12 11:20		Reporting		Instr Dil'n	Prep		Analysis			Reference
	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time		
Antimony	1.2	0.4	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Arsenic	9.4	0.6	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Beryllium	0.3	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Cadmium	0.4	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Chromium	18	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Copper	31	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Lead	41	0.6	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Mercury	< 0.18	0.18	ug/g	1	SAS	10/12/12	5564	10/12/12	17:07	SW7471B	
Nickel	22	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	
Zinc	82	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:33	SW3051A6010C	

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-021

Sample ID: CA-SS-2

Matrix: Solid

Percent Dry: 91.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:10

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	1.1	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Arsenic	6.8	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Beryllium	0.3	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Cadmium	0.5	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Chromium	20	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Copper	16	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Lead	30	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Mercury	< 0.16	0.16	ug/g	1	SAS	10/12/12	5564	10/12/12	17:09	SW7471B
Nickel	12	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Selenium	< 2	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Silver	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C
Zinc	80	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:40	SW3051A6010C

Sample#: 25123-022

Sample ID: CA-SS-3

Matrix: Solid

Percent Dry: 93.4% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:15

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	15	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Arsenic	5.2	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Beryllium	0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Cadmium	4.4	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Chromium	13	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Copper	47	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Lead	190	0.5	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5564	10/12/12	17:11	SW7471B
Nickel	14	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Selenium	< 2	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Silver	0.8	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C
Zinc	310	2	ug/g	1	SAS	10/15/12	5571	10/15/12	23:48	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-023

Sample ID: CA-SS-4

Matrix: Solid

Percent Dry: 91.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 12:45

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	3.6	0.4	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Arsenic	15	0.6	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Beryllium	0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Cadmium	1.5	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Chromium	20	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Copper	32	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Lead	74	0.6	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Mercury	0.48	0.16	ug/g	1	SAS	10/12/12	5564	10/12/12	17:13	SW7471B
Nickel	22	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Thallium	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C
Zinc	140	3	ug/g	1	SAS	10/15/12	5571	10/15/12	23:55	SW3051A6010C

Sample#: 25123-024

Sample ID: CA-SS-5

Matrix: Solid

Percent Dry: 81.2% Results expressed on a dry weight basis.

Sampled: 10/3/12 13:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	66	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Arsenic	26	0.7	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Beryllium	0.4	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Cadmium	21	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Chromium	65	4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Copper	160	4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Lead	1500	0.7	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Mercury	0.36	0.18	ug/g	1	SAS	10/12/12	5565	10/12/12	18:43	SW7471B
Nickel	72	4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Selenium	< 4	4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Silver	4.4	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Thallium	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:02	SW3051A6010C
Zinc	6100	35	ug/g	10	SAS	10/15/12	5571	10/19/12	18:28	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-025

Sample ID: CA-CC-1

Matrix: Solid

Percent Dry: 96.4% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:30

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	2.4	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Arsenic	3.5	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Beryllium	0.7	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Cadmium	0.4	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Chromium	23	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Copper	68	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Lead	84	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Mercury	0.15	0.15	ug/g	1	SAS	10/12/12	5565	10/12/12	17:48	SW7471B
Nickel	17	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Silver	1.6	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C
Zinc	170	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:09	SW3051A6010C

Sample#: 25123-026

Sample ID: CA-CC-2

Matrix: Solid

Percent Dry: 96.9% Results expressed on a dry weight basis.

Sampled: 10/3/12 9:58

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	1.5	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Arsenic	3.2	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Beryllium	0.4	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Cadmium	0.3	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Chromium	20	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Copper	37	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Lead	13	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Mercury	0.37	0.15	ug/g	1	SAS	10/12/12	5565	10/12/12	17:50	SW7471B
Nickel	11	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C
Zinc	92	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:17	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-027

Sample ID: CA-CC-3

Matrix: Solid

Percent Dry: 96.5% Results expressed on a dry weight basis.

Sampled: 10/3/12 10:40

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	0.9	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Arsenic	5.5	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Beryllium	0.6	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Cadmium	0.4	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Chromium	25	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Copper	45	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Lead	34	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5565	10/12/12	17:52	SW7471B
Nickel	18	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C
Zinc	83	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:24	SW3051A6010C

Sample#: 25123-028

Sample ID: CA-CC-4

Matrix: Solid

Percent Dry: 96.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:25

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	4.7	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Arsenic	5.4	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Beryllium	0.7	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Cadmium	2.1	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Chromium	22	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Copper	62	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Lead	76	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5565	10/12/12	17:54	SW7471B
Nickel	18	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C
Zinc	130	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:31	SW3051A6010C

Project ID: MERC 12001150

Job ID: 25123

Sample#: 25123-029

Sample ID: CA-CC-5

Matrix: Solid

Percent Dry: 96.7% Results expressed on a dry weight basis.

Sampled: 10/3/12 11:48

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	13	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Arsenic	6.0	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Beryllium	0.6	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Cadmium	6.9	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Chromium	24	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Copper	77	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Lead	130	0.6	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Mercury	< 0.15	0.15	ug/g	1	SAS	10/12/12	5565	10/12/12	17:55	SW7471B
Nickel	23	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Selenium	< 3	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Silver	< 0.4	0.4	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Thallium	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C
Zinc	260	3	ug/g	1	SAS	10/15/12	5571	10/16/12	0:39	SW3051A6010C

Sample#: 25123-030

Sample ID: STACK

Matrix: Solid

Percent Dry: 99.8% Results expressed on a dry weight basis.

Sampled: 10/3/12 14:50

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Antimony	1.8	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Arsenic	< 0.5	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Beryllium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Cadmium	0.3	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Chromium	4	2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Copper	30	2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Lead	18	0.5	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Mercury	< 0.14	0.14	ug/g	1	SAS	10/12/12	5565	10/12/12	17:57	SW7471B
Nickel	17	2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Selenium	< 2	2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Silver	< 0.3	0.3	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Thallium	< 0.2	0.2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:01	SW3051A6010C
Zinc	32	2	ug/g	1	SAS	10/15/12	5571	10/16/12	1:08	SW3051A6010C

Quality Control Report



124 Heritage Avenue Unit 16
Portsmouth, NH 03801
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Case Narrative

Lab # 25123

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 0 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

ABN: Sample CA-CC-1 and CA-CC-3 did not meet acceptance criteria for the surrogate 2-fluorophenol and 2,4,6-tribromophenol as a result of interferences caused by large co-eluting nontarget compounds.

ABN: Sample CA-CC-2 did not meet acceptance criteria for the surrogate 2-fluorophenol, phenol-D5 and 2,4,6-tribromophenol as a result of interferences caused by large co-eluting nontarget compounds.

PCB: The surrogates were diluted out of the calibration range in the following sample: CA-SB-5 5-10

Laboratory Control Sample Results

ABN: The LCS/D5556 did not meet the acceptance criteria for hexachlorocyclopentadiene. Since <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

VOC: The LCS5548 did not meet the acceptance criteria for bromomethane. The LCSD showed acceptable recovery. This compound is known to be problematic in the method. Since <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

EPH: The LCS/D5547 showed acceptable percent recovery for all analytes, however the RPD between the LCS and LCSD was above the limit for 5 analytes.

EPH: The LCS/D5572 showed acceptable percent recovery for all analytes, however the RPD between the LCS and LCSD was above the limit for fluoranthene.

PCB: The LCS/D5551 showed acceptable percent recovery, however the RPD between the LCS and LCSD was above the limit for both PCB-1016 and PCB-1260.

PCB: The LCS/D5561 showed acceptable percent recovery, however the RPD between the LCS and LCSD was above the limit for both PCB-1016 and PCB-1260.

Thallium: The CRM/D5571 showed acceptable recovery, however the RPD between the CRM and CRMD was above the established limit.



Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Metals: The percent recovery for lead in the lab matrix spike of CA-SB-2 0-1 was 152%, outside the acceptance criteria of 75-125%. This is caused by the high native sample concentration of this analyte relative to the spiked concentration.

Other

EPH: The surrogates in sample CA-SB-12 15-18, CA-SB-13 10-12 and SB-DUP-1 showed surrogate issues as a result of hydrocarbons present in the sample. The chromatograms are included for reference.

EPH: In sample CA-SB-3 10-12.5 no ranges were detected above the method reporting limit; therefore fractionation was not required.

EPH: The surrogate showed recovery outside the acceptance limits for sample CA-SB-2 0-1. Results from re-extraction did not compare well with the original extraction. Both sets of data are provided, the replicate reported as 25123-032.

PAH: Dilution was required due to hydrocarbon interferences for the following samples: CA-MW-8.

PCB: Sample dilution was required for CA-CC-1 and CA-CC-2 due to matrix interferences.

PCB: Sample CA-CC-5 showed elevated reporting limit due to matrix interference. Its surrogate is above the acceptance criteria. Since no targets were detected above the quantitation limit, there is no impact to the data.

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Association Table -

Analysis		QC Number	Field ID	Lab ID
EPH in solids by	MA EPH			
		5547	CA-SB-2 0-1 CA-SB-3 10-12.5 CA-SB-6 12.5-15 CA-SB-7 7.5-10 CA-SB-11 10-15 DUP-2 CA-SB-12 15-18 CA-SB-13 10-12 SB-DUP-1 CA-SB-14 0-3 CA-SB-15 5-5.5	25123-002 25123-003 25123-007 25123-008 25123-013 25123-014 25123-015 25123-016 25123-017 25123-018 25123-019
EPH in solids by	MA EPH			
		5572	CA-SB-4 0-2 CA-SB-2 0-1 Re-Extraction	25123-004 25123-032
VPH in solids by MA	MA VPH			
		5549	CA-SB-2.5-5 CA-SB-4 0-2 CA-SB-6 12.5-15 CA-SB-7 7.5-10 CA-SB-11 10-15 DUP-2 CA-SB-14 0-3 CA-SB-15 5-5.5	25123-001 25123-004 25123-007 25123-008 25123-013 25123-014 25123-018 25123-019
Antimony in solids by	SW3051A6010C			
		5553	CA-SB-2.5-5 CA-SB-2 0-1 CA-SB-3 10-12.5 CA-SB-4 0-2 CA-SB-5 5-10 CA-SB-6 0-5 CA-SB-6 12.5-15 CA-SB-7 7.5-10 CA-SB-8 0-2.5 CA-SB-8 5-10 CA-SB-10 10-15 CA-SB-11 10-15 CA-SB-12 15-18	25123-001 25123-002 25123-003 25123-004 25123-005 25123-006 25123-007 25123-008 25123-009 25123-010 25123-012 25123-013 25123-015
Antimony in solids by	SW3051A6010C			
		5571	CA-SB-13 10-12 SB-DUP-1 CA-SS-1 CA-SS-2 CA-SS-3 CA-SS-4 CA-SS-5 CA-CC-1 CA-CC-2 CA-CC-3 CA-CC-4	25123-016 25123-017 25123-020 25123-021 25123-022 25123-023 25123-024 25123-025 25123-026 25123-027 25123-028

Analysis	QC Number	Field ID	Lab ID
	5571	CA-SS-5	25123-029
		STACK	25123-030
PCBs in soil by 8082	SW3540C8082A		
	5551	CA-SB-3 10-12.5	25123-003
		CA-SB-5 5-10	25123-005
		CA-SB-6 0-5	25123-006
		CA-SB-6 12.5-15	25123-007
		CA-SB-7 7.5-10	25123-008
		CA-SB-8 0-2.5	25123-009
		CA-SB-8 5-10	25123-010
		CA-SB-9 0-4	25123-011
		CA-SB-11 10-15	25123-013
		CA-SB-12 15-18	25123-015
		CA-SB-13 10-12	25123-016
		SB-DUP-1	25123-017
		CA-SS-1	25123-020
		CA-SS-2	25123-021
		CA-SS-3	25123-022
PCBs in soil by 8082	SW3540C8082A		
	5561	CA-SB-10 10-15	25123-012
		CA-SS-4	25123-023
		CA-SS-5	25123-024
		CA-CC-3	25123-027
		CA-CC-4	25123-028
		STACK	25123-030
Acid & Base/Neutral	SW3546/8270D		
	5556	CA-SB-4 0-2	25123-004
		CA-SB-5 5-10	25123-005
		CA-SB-6 12.5-15	25123-007
		CA-SB-7 7.5-10	25123-008
		CA-SB-10 10-15	25123-012
		CA-SB-11 10-15	25123-013
		DUP-2	25123-014
		CA-CC-1	25123-025
		CA-CC-2	25123-026
		CA-CC-3	25123-027
PAHs in solid by 8270	SW3550B8270D		
	5555	CA-SB-6 0-5	25123-006
		CA-SB-8 0-2.5	25123-009
		CA-SB-8 5-10	25123-010
		CA-SS-1	25123-020
		CA-SS-2	25123-021
		CA-SS-3	25123-022
VOCs in solids by 8260	SW5035A8260B		
	5548	CA-SB-4 0-2	25123-004
		CA-SB-5 5-10	25123-005
		CA-SB-6 12.5-15	25123-007
		CA-SB-7 7.5-10	25123-008
		CA-SB-10 10-15	25123-012
		CA-SB-11 10-15	25123-013
		DUP-2	25123-014
		CA-CC-1	25123-025

Analysis	QC Number	Field ID	Lab ID
	5548	CA-CC-2	25123-026
		CA-CC-3	25123-027
		Trip Blank	25123-031
Mercury in solids by	SW7471B		
	5564	CA-SB-2 0-1	25123-002
		CA-SB-3 10-12.5	25123-003
		CA-SB-4 0-2	25123-004
		CA-SB-5 5-10	25123-005
		CA-SB-6 0-5	25123-006
		CA-SB-6 12.5-15	25123-007
		CA-SB-7 7.5-10	25123-008
		CA-SB-8 0-2.5	25123-009
		CA-SB-8 5-10	25123-010
		CA-SB-10 10-15	25123-012
		CA-SB-11 10-15	25123-013
		CA-SB-12 15-18	25123-015
		CA-SB-13 10-12	25123-016
		SB-DUP-1	25123-017
		CA-SS-1	25123-020
		CA-SS-2	25123-021
		CA-SS-3	25123-022
		CA-SS-4	25123-023
Mercury in solids by	SW7471B		
	5565	CA-SS-5	25123-024
		CA-CC-1	25123-025
		CA-CC-2	25123-026
		CA-CC-3	25123-027
		CA-CC-4	25123-028
		CA-SS-5	25123-029
		STACK	25123-030

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA VPH	MB5549	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		<	5	ug/g				
		Unadjusted C9-C12 Aliphatics		<	5	ug/g				
		methyl t-butyl ether (MTBE)		<	0.1	ug/g				
		benzene		<	0.1	ug/g				
		toluene		<	0.1	ug/g				
		ethylbenzene		<	0.1	ug/g				
		m&p xylenes		<	0.1	ug/g				
		o-xylene		<	0.1	ug/g				
		naphthalene		<	0.2	ug/g				
		C9-C10 Aromatics		<	5	ug/g				
		2,5-dibromotoluene as		102	%			70	130	
		2,5-dibromotoluene as		105	%			70	130	
		a,a,a-trifluorotoluene SUR		104	%			70	130	
MA VPH	MLCS5549	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		7.7	ug/g	7.5	103	70	130	
		Unadjusted C9-C12 Aliphatics		5.0	ug/g	5.5	91	70	130	
		methyl t-butyl ether (MTBE)		1.5	ug/g	1.5	97	70	130	
		benzene		0.5	ug/g	0.5	91	70	130	
		toluene		1.4	ug/g	1.5	94	70	130	
		ethylbenzene		0.4	ug/g	0.5	85	70	130	
		m&p xylenes		1.9	ug/g	2	96	70	130	
		o-xylene		1.0	ug/g	1	95	70	130	
		naphthalene		1.0	ug/g	1	98	70	130	
		C9-C10 Aromatics		1.0	ug/g	1	103	70	130	
		2,5-dibromotoluene as		106	%			70	130	
		2,5-dibromotoluene as		109	%			70	130	
		a,a,a-trifluorotoluene SUR		111	%			70	130	
MA VPH	MLCSD5549	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		7.7	ug/g	7.5	103	70	130	0
		Unadjusted C9-C12 Aliphatics		5.0	ug/g	5.5	91	70	130	1
		methyl t-butyl ether (MTBE)		1.4	ug/g	1.5	96	70	130	1
		benzene		0.5	ug/g	0.5	90	70	130	0
		toluene		1.4	ug/g	1.5	93	70	130	1
		ethylbenzene		0.4	ug/g	0.5	80	70	130	6
		m&p xylenes		1.9	ug/g	2	97	70	130	1
		o-xylene		1.0	ug/g	1	96	70	130	1
		naphthalene		0.9	ug/g	1	92	70	130	5
		C9-C10 Aromatics		1.0	ug/g	1	102	70	130	1
		2,5-dibromotoluene as		103	%			70	130	
		2,5-dibromotoluene as		107	%			70	130	
		a,a,a-trifluorotoluene SUR		107	%			70	130	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB5548	dichlorodifluoromethane		<	0.1	ug/g				
		chloromethane		<	0.1	ug/g				
		vinyl chloride		<	0.1	ug/g				
		bromomethane		<	0.2	ug/g				
		chloroethane		<	0.1	ug/g				
		trichlorofluoromethane		<	0.1	ug/g				
		diethyl ether		<	0.5	ug/g				
		acetone		<	2.5	ug/g				
		1,1-dichloroethene		<	0.1	ug/g				
		methylene chloride		<	0.2	ug/g				
		carbon disulfide		<	0.1	ug/g				
		methyl t-butyl ether (MTBE)		<	0.1	ug/g				
		trans-1,2-dichloroethene		<	0.1	ug/g				
		1,1-dichloroethane		<	0.1	ug/g				
		2-butanone (MEK)		<	0.5	ug/g				
		2,2-dichloropropane		<	0.1	ug/g				
		cis-1,2-dichloroethene		<	0.1	ug/g				
		chloroform		<	0.1	ug/g				
		bromochloromethane		<	0.1	ug/g				
		tetrahydrofuran (THF)		<	0.5	ug/g				
		1,1,1-trichloroethane		<	0.1	ug/g				
		1,1-dichloropropene		<	0.1	ug/g				
		carbon tetrachloride		<	0.1	ug/g				
		1,2-dichloroethane		<	0.1	ug/g				
		benzene		<	0.1	ug/g				
		trichloroethene		<	0.1	ug/g				
		1,2-dichloropropane		<	0.1	ug/g				
		bromodichloromethane		<	0.1	ug/g				
		dibromomethane		<	0.1	ug/g				
		4-methyl-2-pentanone (MIBK)		<	0.5	ug/g				
		cis-1,3-dichloropropene		<	0.1	ug/g				
		toluene		<	0.1	ug/g				
		trans-1,3-dichloropropene		<	0.1	ug/g				
		2-hexanone		<	0.5	ug/g				
		1,1,2-trichloroethane		<	0.1	ug/g				
		1,3-dichloropropane		<	0.1	ug/g				
		tetrachloroethene		<	0.1	ug/g				
		dibromochloromethane		<	0.1	ug/g				
		1,2-dibromoethane (EDB)		<	0.1	ug/g				
		chlorobenzene		<	0.1	ug/g				
		1,1,1,2-tetrachloroethane		<	0.1	ug/g				
		ethylbenzene		<	0.1	ug/g				
		m&p-xylenes		<	0.1	ug/g				
		o-xylene		<	0.1	ug/g				
		styrene		<	0.1	ug/g				
		bromoform		<	0.1	ug/g				
		isopropylbenzene		<	0.1	ug/g				
		1,1,2,2-tetrachloroethane		<	0.1	ug/g				
		1,2,3-trichloropropane		<	0.1	ug/g				
		n-propylbenzene		<	0.1	ug/g				
		bromobenzene		<	0.1	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MB5548	1,3,5-trimethylbenzene		<	0.1	ug/g				
		2-chlorotoluene		<	0.1	ug/g				
		4-chlorotoluene		<	0.1	ug/g				
		tert-butylbenzene		<	0.1	ug/g				
		1,2,4-trimethylbenzene		<	0.1	ug/g				
		sec-butylbenzene		<	0.1	ug/g				
		1,3-dichlorobenzene		<	0.1	ug/g				
		4-isopropyltoluene		<	0.1	ug/g				
		1,4-dichlorobenzene		<	0.1	ug/g				
		1,2-dichlorobenzene		<	0.1	ug/g				
		n-butylbenzene		<	0.1	ug/g				
		1,2-dibromo-3-chloropropane		<	0.1	ug/g				
		1,2,4-trichlorobenzene		<	0.1	ug/g				
		hexachlorobutadiene		<	0.1	ug/g				
		naphthalene		<	0.2	ug/g				
		1,2,3-trichlorobenzene		<	0.1	ug/g				
		dibromofluoromethane SUR		99	%		78	114		
		toluene-D8 SUR		97	%		88	110		
		4-bromofluorobenzene SUR		104	%		86	115		
		a,a,a-trifluorotoluene SUR		99	%		70	130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS5548	dichlorodifluoromethane		0.8	ug/g	1	80	70 130		
		chloromethane		0.8	ug/g	1	82	70 130		
		vinyl chloride		0.9	ug/g	1	85	70 130		
		bromomethane		0.6	ug/g	1	61 *	70 130		
		chloroethane		1.0	ug/g	1	95	70 130		
		trichlorofluoromethane		0.9	ug/g	1	91	70 130		
		diethyl ether		0.9	ug/g	1	95	70 130		
		acetone	<	2.5	ug/g	1	92			
		1,1-dichloroethene		0.8	ug/g	1	83	70 130		
		methylene chloride		0.9	ug/g	1	95	70 130		
		carbon disulfide		1.0	ug/g	1	98	70 130		
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	99	70 130		
		trans-1,2-dichloroethene		0.9	ug/g	1	92	70 130		
		1,1-dichloroethane		0.9	ug/g	1	91	70 130		
		2-butanone (MEK)		1.1	ug/g	1	105	70 130		
		2,2-dichloropropane		0.9	ug/g	1	90	70 130		
		cis-1,2-dichloroethene		1.0	ug/g	1	95	70 130		
		chloroform		1.0	ug/g	1	97	70 130		
		bromochloromethane		0.9	ug/g	1	91	70 130		
		tetrahydrofuran (THF)		0.9	ug/g	1	89	70 130		
		1,1,1-trichloroethane		0.9	ug/g	1	93	70 130		
		1,1-dichloropropene		1.0	ug/g	1	95	70 130		
		carbon tetrachloride		0.9	ug/g	1	90	70 130		
		1,2-dichloroethane		0.9	ug/g	1	93	70 130		
		benzene		0.9	ug/g	1	92	70 130		
		trichloroethene		1.0	ug/g	1	95	70 130		
		1,2-dichloropropane		0.9	ug/g	1	92	70 130		
		bromodichloromethane		0.9	ug/g	1	94	70 130		
		dibromomethane		0.9	ug/g	1	94	70 130		
		4-methyl-2-pentanone (MIBK)		1.0	ug/g	1	105	70 130		
		cis-1,3-dichloropropene		0.9	ug/g	1	93	70 130		
		toluene		0.9	ug/g	1	95	70 130		
		trans-1,3-dichloropropene		0.9	ug/g	1	89	70 130		
		2-hexanone		1.1	ug/g	1	110	70 130		
		1,1,2-trichloroethane		0.9	ug/g	1	92	70 130		
		1,3-dichloropropane		0.9	ug/g	1	92	70 130		
		tetrachloroethene		0.9	ug/g	1	90	70 130		
		dibromochloromethane		0.9	ug/g	1	89	70 130		
		1,2-dibromoethane (EDB)		0.9	ug/g	1	92	70 130		
		chlorobenzene		0.9	ug/g	1	93	70 130		
		1,1,1,2-tetrachloroethane		0.9	ug/g	1	93	70 130		
		ethylbenzene		0.9	ug/g	1	93	70 130		
		m&p-xlenes		2.0	ug/g	2	98	70 130		
		o-xylene		1.0	ug/g	1	97	70 130		
		styrene		1.0	ug/g	1	101	70 130		
		bromoform		1.0	ug/g	1	97	70 130		
		isopropylbenzene		0.9	ug/g	1	94	70 130		
		1,1,2,2-tetrachloroethane		1.0	ug/g	1	99	70 130		
		1,2,3-trichloropropane		0.9	ug/g	1	94	70 130		
		n-propylbenzene		0.9	ug/g	1	90	70 130		
		bromobenzene		0.9	ug/g	1	93	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCS5548	1,3,5-trimethylbenzene		0.9	ug/g	1	91	70 130		
		2-chlorotoluene		0.9	ug/g	1	93	70 130		
		4-chlorotoluene		1.0	ug/g	1	98	70 130		
		tert-butylbenzene		0.9	ug/g	1	90	70 130		
		1,2,4-trimethylbenzene		0.9	ug/g	1	92	70 130		
		sec-butylbenzene		1.0	ug/g	1	96	70 130		
		1,3-dichlorobenzene		1.0	ug/g	1	97	70 130		
		4-isopropyltoluene		1.0	ug/g	1	95	70 130		
		1,4-dichlorobenzene		1.0	ug/g	1	96	70 130		
		1,2-dichlorobenzene		1.0	ug/g	1	98	70 130		
		n-butylbenzene		1.0	ug/g	1	96	70 130		
		1,2-dibromo-3-chloropropane		1.0	ug/g	1	99	70 130		
		1,2,4-trichlorobenzene		0.9	ug/g	1	91	70 130		
		hexachlorobutadiene		0.9	ug/g	1	93	70 130		
		naphthalene		1.0	ug/g	1	100	70 130		
		1,2,3-trichlorobenzene		1.0	ug/g	1	97	70 130		
		dibromofluoromethane SUR		102	%			78 114		
		toluene-D8 SUR		100	%			88 110		
		4-bromofluorobenzene SUR		110	%			86 115		
		a,a,a-trifluorotoluene SUR		102	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD5548	dichlorodifluoromethane		0.9	ug/g	1	92	70 130	14	30
		chloromethane		0.9	ug/g	1	92	70 130	11	30
		vinyl chloride		1.0	ug/g	1	98	70 130	13	30
		bromomethane		0.7	ug/g	1	73	70 130	18	30
		chloroethane		1.1	ug/g	1	108	70 130	12	30
		trichlorofluoromethane		1.0	ug/g	1	101	70 130	11	30
		diethyl ether		1.0	ug/g	1	103	70 130	8	30
		acetone	<	2.5	ug/g	1	95		3	30
		1,1-dichloroethene		0.9	ug/g	1	90	70 130	8	30
		methylene chloride		1.1	ug/g	1	107	70 130	12	30
		carbon disulfide		1.1	ug/g	1	108	70 130	10	30
		methyl t-butyl ether (MTBE)		1.0	ug/g	1	99	70 130	0	30
		trans-1,2-dichloroethene		1.0	ug/g	1	99	70 130	7	30
		1,1-dichloroethane		1.0	ug/g	1	99	70 130	8	30
		2-butanone (MEK)		1.0	ug/g	1	102	70 130	3	30
		2,2-dichloropropane		1.0	ug/g	1	96	70 130	6	30
		cis-1,2-dichloroethene		1.1	ug/g	1	105	70 130	10	30
		chloroform		1.0	ug/g	1	104	70 130	7	30
		bromochloromethane		1.0	ug/g	1	102	70 130	11	30
		tetrahydrofuran (THF)		0.8	ug/g	1	80	70 130	10	30
		1,1,1-trichloroethane		1.0	ug/g	1	100	70 130	7	30
		1,1-dichloropropene		1.0	ug/g	1	102	70 130	7	30
		carbon tetrachloride		1.0	ug/g	1	97	70 130	7	30
		1,2-dichloroethane		1.0	ug/g	1	99	70 130	6	30
		benzene		1.0	ug/g	1	99	70 130	7	30
		trichloroethene		1.0	ug/g	1	100	70 130	5	30
		1,2-dichloropropane		1.0	ug/g	1	99	70 130	8	30
		bromodichloromethane		1.0	ug/g	1	99	70 130	6	30
		dibromomethane		1.0	ug/g	1	97	70 130	3	30
		4-methyl-2-pentanone (MIBK)		1.0	ug/g	1	103	70 130	2	30
		cis-1,3-dichloropropene		1.0	ug/g	1	102	70 130	9	30
		toluene		1.0	ug/g	1	104	70 130	9	30
		trans-1,3-dichloropropene		1.0	ug/g	1	97	70 130	9	30
		2-hexanone		1.0	ug/g	1	104	70 130	5	30
		1,1,2-trichloroethane		1.0	ug/g	1	100	70 130	8	30
		1,3-dichloropropane		1.0	ug/g	1	98	70 130	6	30
		tetrachloroethene		0.9	ug/g	1	94	70 130	4	30
		dibromochloromethane		1.0	ug/g	1	96	70 130	8	30
		1,2-dibromoethane (EDB)		0.9	ug/g	1	95	70 130	3	30
		chlorobenzene		1.0	ug/g	1	101	70 130	9	30
		1,1,1,2-tetrachloroethane		1.0	ug/g	1	101	70 130	8	30
		ethylbenzene		1.0	ug/g	1	101	70 130	8	30
		m&p-xlenes		2.1	ug/g	2	107	70 130	8	30
		o-xylene		1.1	ug/g	1	107	70 130	9	30
		styrene		1.1	ug/g	1	111	70 130	9	30
		bromoform		1.0	ug/g	1	101	70 130	4	30
		isopropylbenzene		1.0	ug/g	1	102	70 130	9	30
		1,1,2,2-tetrachloroethane		1.0	ug/g	1	97	70 130	2	30
		1,2,3-trichloropropane		0.9	ug/g	1	92	70 130	3	30
		n-propylbenzene		1.0	ug/g	1	97	70 130	8	30
		bromobenzene		1.0	ug/g	1	100	70 130	7	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5035A8260B	MLCSD5548	1,3,5-trimethylbenzene		1.0	ug/g	1	98	70 130	8	30
		2-chlorotoluene		1.0	ug/g	1	98	70 130	6	30
		4-chlorotoluene		1.0	ug/g	1	105	70 130	7	30
		tert-butylbenzene		1.0	ug/g	1	102	70 130	12	30
		1,2,4-trimethylbenzene		1.0	ug/g	1	99	70 130	7	30
		sec-butylbenzene		1.0	ug/g	1	104	70 130	8	30
		1,3-dichlorobenzene		1.0	ug/g	1	104	70 130	7	30
		4-isopropyltoluene		1.0	ug/g	1	101	70 130	6	30
		1,4-dichlorobenzene		1.0	ug/g	1	104	70 130	8	30
		1,2-dichlorobenzene		1.1	ug/g	1	106	70 130	7	30
		n-butylbenzene		1.1	ug/g	1	105	70 130	9	30
		1,2-dibromo-3-chloropropane		1.0	ug/g	1	96	70 130	3	30
		1,2,4-trichlorobenzene		1.0	ug/g	1	97	70 130	7	30
		hexachlorobutadiene		1.0	ug/g	1	101	70 130	8	30
		naphthalene		1.0	ug/g	1	100	70 130	1	30
		1,2,3-trichlorobenzene		1.0	ug/g	1	101	70 130	4	30
		dibromofluoromethane SUR		98	%			78 114		
		toluene-D8 SUR		102	%			88 110		
		4-bromofluorobenzene SUR		108	%			86 115		
		a,a,a-trifluorotoluene SUR		104	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK5547	naphthalene		<	0.1	ug/g				
		2-methylnaphthalene		<	0.1	ug/g				
		phenanthrene		<	0.1	ug/g				
		acenaphthene		<	0.1	ug/g				
		acenaphthylene		<	0.1	ug/g				
		fluorene		<	0.1	ug/g				
		anthracene		<	0.1	ug/g				
		fluoranthene		<	0.1	ug/g				
		pyrene		<	0.1	ug/g				
		benzo(a)anthracene		<	0.1	ug/g				
		chrysene		<	0.1	ug/g				
		benzo(b)fluoranthene		<	0.1	ug/g				
		benzo(k)fluoranthene		<	0.1	ug/g				
		benzo(a)pyrene		<	0.1	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.1	ug/g				
		dibenzo(a,h)anthracene		<	0.1	ug/g				
		benzo(g,h,i)perylene		<	0.1	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			62	%		40	140	
		o-terphenyl SUR			67	%		40	140	
		2-fluorobiphenyl SUR			105	%		40	140	
		2-bromonaphthalene SUR			104	%		40	140	
MA EPH	LCS5547	naphthalene		2.6	ug/g	6	43	40	140	
		2-methylnaphthalene		2.6	ug/g	6	44	40	140	
		phenanthrene		3.1	ug/g	6	51	40	140	
		acenaphthene		2.8	ug/g	6	47	40	140	
		acenaphthylene		2.9	ug/g	6	48	40	140	
		fluorene		2.8	ug/g	6	47	40	140	
		anthracene		2.9	ug/g	6	48	40	140	
		fluoranthene		3.0	ug/g	6	51	40	140	
		pyrene		3.5	ug/g	6	59	40	140	
		benzo(a)anthracene		3.5	ug/g	6	58	40	140	
		chrysene		3.4	ug/g	6	58	40	140	
		benzo(b)fluoranthene		3.7	ug/g	6	62	40	140	
		benzo(k)fluoranthene		3.5	ug/g	6	59	40	140	
		benzo(a)pyrene		3.4	ug/g	6	57	40	140	
		indeno(1,2,3-cd)pyrene		3.5	ug/g	6	59	40	140	
		dibenzo(a,h)anthracene		3.7	ug/g	6	61	40	140	
		benzo(g,h,i)perylene		3.7	ug/g	6	62	40	140	
		Unadjusted C11-C22 Aromatics			43	ug/g	102	42	40	140
		C9-C18 Aliphatics			25	ug/g	36	68	40	140
		C19-C36 Aliphatics			27	ug/g	48	57	40	140
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			45	%		40	140	
		o-terphenyl SUR			47	%		40	140	
		2-fluorobiphenyl SUR			97	%		40	140	
		2-bromonaphthalene SUR			82	%		40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
MA EPH	LCSD5547	naphthalene		3.4	ug/g	6	57	40 140	27	*	25
		2-methylnaphthalene		3.4	ug/g	6	56	40 140	24		25
		phenanthrene		3.7	ug/g	6	62	40 140	19		25
		acenaphthene		3.5	ug/g	6	59	40 140	23		25
		acenaphthylene		3.6	ug/g	6	60	40 140	23		25
		fluorene		3.7	ug/g	6	61	40 140	26	*	25
		anthracene		3.8	ug/g	6	63	40 140	27	*	25
		fluoranthene		3.5	ug/g	6	58	40 140	13		25
		pyrene		4.3	ug/g	6	71	40 140	19		25
		benzo(a)anthracene		4.0	ug/g	6	66	40 140	13		25
		chrysene		4.1	ug/g	6	68	40 140	17		25
		benzo(b)fluoranthene		3.7	ug/g	6	62	40 140	0		25
		benzo(k)fluoranthene		4.8	ug/g	6	79	40 140	30	*	25
		benzo(a)pyrene		4.0	ug/g	6	66	40 140	15		25
		indeno(1,2,3-cd)pyrene		4.0	ug/g	6	67	40 140	12		25
		dibenzo(a,h)anthracene		4.0	ug/g	6	66	40 140	9		25
		benzo(g,h,i)perylene		3.9	ug/g	6	65	40 140	4		25
		Unadjusted C11-C22 Aromatics		53	ug/g	102	52	40 140	21		25
		C9-C18 Aliphatics		30	ug/g	36	83	40 140	19		25
		C19-C36 Aliphatics		38	ug/g	48	79	40 140	32	*	25
		C11-C22 Aromatics	<	20	ug/g						
		1-chloro-octadecane SUR		59	%			40 140			
		o-terphenyl SUR		65	%			40 140			
		2-fluorobiphenyl SUR		103	%			40 140			
		2-bromonaphthalene SUR		86	%			40 140			

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK5572	naphthalene		<	0.1	ug/g				
		2-methylnaphthalene		<	0.1	ug/g				
		phenanthrene		<	0.1	ug/g				
		acenaphthene		<	0.1	ug/g				
		acenaphthylene		<	0.1	ug/g				
		fluorene		<	0.1	ug/g				
		anthracene		<	0.1	ug/g				
		fluoranthene		<	0.1	ug/g				
		pyrene		<	0.1	ug/g				
		benzo(a)anthracene		<	0.1	ug/g				
		chrysene		<	0.1	ug/g				
		benzo(b)fluoranthene		<	0.1	ug/g				
		benzo(k)fluoranthene		<	0.1	ug/g				
		benzo(a)pyrene		<	0.1	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.1	ug/g				
		dibenzo(a,h)anthracene		<	0.1	ug/g				
		benzo(g,h,i)perylene		<	0.1	ug/g				
		Unadjusted C11-C22 Aromatics		<	20	ug/g				
		C9-C18 Aliphatics		<	20	ug/g				
		C19-C36 Aliphatics		<	20	ug/g				
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			62	%		40	140	
		o-terphenyl SUR			69	%		40	140	
		2-fluorobiphenyl SUR			99	%		40	140	
		2-bromonaphthalene SUR			96	%		40	140	
MA EPH	LCS5572	naphthalene		3.8	ug/g	6	63	40	140	
		2-methylnaphthalene		3.7	ug/g	6	62	40	140	
		phenanthrene		4.4	ug/g	6	74	40	140	
		acenaphthene		4.0	ug/g	6	67	40	140	
		acenaphthylene		4.3	ug/g	6	71	40	140	
		fluorene		3.6	ug/g	6	59	40	140	
		anthracene		4.3	ug/g	6	72	40	140	
		fluoranthene		4.5	ug/g	6	74	40	140	
		pyrene		5.2	ug/g	6	87	40	140	
		benzo(a)anthracene		4.9	ug/g	6	81	40	140	
		chrysene		5.0	ug/g	6	83	40	140	
		benzo(b)fluoranthene		4.8	ug/g	6	79	40	140	
		benzo(k)fluoranthene		5.3	ug/g	6	89	40	140	
		benzo(a)pyrene		4.8	ug/g	6	80	40	140	
		indeno(1,2,3-cd)pyrene		4.9	ug/g	6	82	40	140	
		dibenzo(a,h)anthracene		5.0	ug/g	6	83	40	140	
		benzo(g,h,i)perylene		4.9	ug/g	6	82	40	140	
		Unadjusted C11-C22 Aromatics			48	ug/g	102	48	40	140
		C9-C18 Aliphatics			32	ug/g	36	89	40	140
		C19-C36 Aliphatics			42	ug/g	48	87	40	140
		C11-C22 Aromatics		<	20	ug/g				
		1-chloro-octadecane SUR			64	%		40	140	
		o-terphenyl SUR			62	%		40	140	
		2-fluorobiphenyl SUR			85	%		40	140	
		2-bromonaphthalene SUR			74	%		40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD5572	naphthalene		3.9	ug/g	6	65	40 140	3	25
		2-methylnaphthalene		3.9	ug/g	6	65	40 140	4	25
		phenanthrene		4.3	ug/g	6	72	40 140	2	25
		acenaphthene		4.1	ug/g	6	69	40 140	3	25
		acenaphthylene		4.2	ug/g	6	71	40 140	1	25
		fluorene		3.9	ug/g	6	66	40 140	10	25
		anthracene		4.0	ug/g	6	67	40 140	6	25
		fluoranthene		3.3	ug/g	6	55	40 140	29 *	25
		pyrene		4.8	ug/g	6	79	40 140	9	25
		benzo(a)anthracene		4.7	ug/g	6	78	40 140	3	25
		chrysene		4.9	ug/g	6	82	40 140	2	25
		benzo(b)fluoranthene		4.3	ug/g	6	72	40 140	10	25
		benzo(k)fluoranthene		5.5	ug/g	6	92	40 140	3	25
		benzo(a)pyrene		4.6	ug/g	6	77	40 140	4	25
		indeno(1,2,3-cd)pyrene		4.7	ug/g	6	78	40 140	5	25
		dibenzo(a,h)anthracene		4.9	ug/g	6	81	40 140	2	25
		benzo(g,h,i)perylene		4.6	ug/g	6	77	40 140	6	25
		Unadjusted C11-C22 Aromatics		49	ug/g	102	48	40 140	1	25
		C9-C18 Aliphatics		31	ug/g	36	85	40 140	5	25
		C19-C36 Aliphatics		38	ug/g	48	78	40 140	11	25
		C11-C22 Aromatics	<	20	ug/g					
		1-chloro-octadecane SUR		57	%			40 140		
		o-terphenyl SUR		59	%			40 140		
		2-fluorobiphenyl SUR		100	%			40 140		
		2-bromonaphthalene SUR		83	%			40 140		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3540C8082A	BLK5551	PCB-1016		<	0.02	ug/g				
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		<	0.02	ug/g				
		tetrachloro-m-xylene SUR		73	%		30	150		
		decachlorobiphenyl SUR		66	%		30	150		
SW3540C8082A	LCS5551	PCB-1016		0.10	ug/g	0.2	50	40	140	
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.10	ug/g	0.2	49	40	140	
		tetrachloro-m-xylene SUR		35	%		30	150		
		decachlorobiphenyl SUR		48	%		30	150		
SW3540C8082A	LCSD5551	PCB-1016		0.19	ug/g	0.2	93	40	140	60 * 30
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.17	ug/g	0.2	87	40	140	55 * 30
		tetrachloro-m-xylene SUR		69	%		30	150		
		decachlorobiphenyl SUR		77	%		30	150		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3540C8082A	BLK5561	PCB-1016		<	0.02	ug/g				
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		<	0.02	ug/g				
		tetrachloro-m-xylene SUR		56	%		30	150		
		decachlorobiphenyl SUR		66	%		30	150		
SW3540C8082A	LCS5561	PCB-1016		0.12	ug/g	0.2	59	40	140	
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.10	ug/g	0.2	52	40	140	
		tetrachloro-m-xylene SUR		45	%		30	150		
		decachlorobiphenyl SUR		58	%		30	150		
SW3540C8082A	LCSD5561	PCB-1016		0.17	ug/g	0.2	86	40	140	36 *
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.15	ug/g	0.2	73	40	140	33 *
		tetrachloro-m-xylene SUR		64	%		30	150		
		decachlorobiphenyl SUR		73	%		30	150		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK5556	N-nitrosodimethylamine		<	0.2	ug/g				
		aniline		<	0.2	ug/g				
		phenol		<	0.2	ug/g				
		2-chlorophenol		<	0.5	ug/g				
		bis(2-chloroethyl)ether		<	0.2	ug/g				
		1,3-dichlorobenzene		<	0.2	ug/g				
		1,4-dichlorobenzene		<	0.2	ug/g				
		1,2-dichlorobenzene		<	0.2	ug/g				
		benzyl alcohol		<	0.2	ug/g				
		2-methylphenol		<	0.2	ug/g				
		bis(2-chloroisopropyl) ether		<	0.2	ug/g				
		hexachloroethane		<	0.2	ug/g				
		N-nitroso-di-N-propylamine		<	0.2	ug/g				
		4-methylphenol		<	0.2	ug/g				
		nitrobenzene		<	0.2	ug/g				
		isophorone		<	0.5	ug/g				
		2-nitrophenol		<	0.2	ug/g				
		2,4-dimethylphenol		<	0.2	ug/g				
		bis(2-chloroethoxy)methane		<	0.5	ug/g				
		2,4-dichlorophenol		<	0.5	ug/g				
		1,2,4-trichlorobenzene		<	0.5	ug/g				
		naphthalene		<	0.05	ug/g				
		benzoic acid		<	5.0	ug/g				
		4-chloroaniline		<	0.2	ug/g				
		hexachlorobutadiene		<	0.2	ug/g				
		4-chloro-3-methylphenol		<	0.2	ug/g				
		2-methylnaphthalene		<	0.05	ug/g				
		hexachlorocyclopentadiene		<	1.0	ug/g				
		2,4,6-trichlorophenol		<	0.2	ug/g				
		2,4,5-trichlorophenol		<	0.2	ug/g				
		2-chloronaphthalene		<	0.5	ug/g				
		2-nitroaniline		<	0.2	ug/g				
		acenaphthylene		<	0.05	ug/g				
		dimethylphthalate		<	0.5	ug/g				
		2,6-dinitrotoluene		<	0.2	ug/g				
		2,4-dinitrotoluene		<	0.2	ug/g				
		acenaphthene		<	0.05	ug/g				
		3-nitroaniline		<	0.2	ug/g				
		2,4-dinitrophenol		<	5.0	ug/g				
		dibenzofuran		<	0.05	ug/g				
		4-nitrophenol		<	1.0	ug/g				
		fluorene		<	0.05	ug/g				
		diethyl phthalate		<	0.5	ug/g				
		4-chlorophenyl phenyl ether		<	0.5	ug/g				
		4-nitroaniline		<	0.5	ug/g				
		4,6-dinitro-2-methylphenol		<	2.0	ug/g				
		azobenzene		<	0.2	ug/g				
		N-nitrosodiphenylamine		<	0.2	ug/g				
		4-bromophenyl phenyl ether		<	0.2	ug/g				
		hexachlorobenzene		<	0.2	ug/g				
		pentachlorophenol		<	1.0	ug/g				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	BLK5556	phenanthrene		<	0.05	ug/g				
		anthracene		<	0.05	ug/g				
		carbazole		<	0.2	ug/g				
		di-n-butylphthalate		<	0.5	ug/g				
		fluoranthene		<	0.05	ug/g				
		benzidine		<	3.0	ug/g				
		pyrene		<	0.05	ug/g				
		butyl benzyl phthalate		<	0.5	ug/g				
		benzo(a)anthracene		<	0.05	ug/g				
		chrysene		<	0.05	ug/g				
		3,3'-dichlorobenzidine		<	3.0	ug/g				
		bis(2-ethylhexyl)phthalate		<	0.5	ug/g				
		di-n-octyl phthalate		<	0.2	ug/g				
		benzo(b)fluoranthene		<	0.05	ug/g				
		benzo(k)fluoranthene		<	0.05	ug/g				
		benzo(a)pyrene		<	0.02	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.05	ug/g				
		dibenzo(a,h)anthracene		<	0.05	ug/g				
		benzo(g,h,i)perylene		<	0.05	ug/g				
		2-fluorophenol SUR		58	%			21	100	
		phenol-D5 SUR		56	%			10	102	
		2,4,6-tribromophenol SUR		59	%			10	123	
		nitrobenzene-D5 SUR		58	%			35	114	
		2-fluorobiphenyl SUR		69	%			43	116	
		p-terphenyl-D14 SUR		78	%			33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS5556	N-nitrosodimethylamine		1.9	ug/g	4	48	40 140		
		aniline		1.9	ug/g	4	46	40 140		
		phenol		2.3	ug/g	4	57	30 130		
		2-chlorophenol		2.5	ug/g	4	63	30 130		
		bis(2-chloroethyl)ether		2.9	ug/g	4	72	40 140		
		1,3-dichlorobenzene		2.2	ug/g	4	55	40 140		
		1,4-dichlorobenzene		2.3	ug/g	4	58	40 140		
		1,2-dichlorobenzene		2.4	ug/g	4	60	40 140		
		benzyl alcohol		2.2	ug/g	4	54	30 130		
		2-methylphenol		2.5	ug/g	4	62	30 130		
		bis(2-chloroisopropyl) ether		2.3	ug/g	4	58	40 140		
		hexachloroethane		2.2	ug/g	4	56	40 140		
		N-nitroso-di-N-propylamine		2.2	ug/g	4	54	40 140		
		4-methylphenol		2.6	ug/g	4	64	30 130		
		nitrobenzene		2.2	ug/g	4	55	40 140		
		isophorone		2.2	ug/g	4	56	40 140		
		2-nitrophenol		2.1	ug/g	4	53	30 130		
		2,4-dimethylphenol		2.6	ug/g	4	65	30 130		
		bis(2-chloroethoxy)methane		2.4	ug/g	4	60	40 140		
		2,4-dichlorophenol		2.7	ug/g	4	67	30 130		
		1,2,4-trichlorobenzene		2.3	ug/g	4	57	40 140		
		naphthalene		2.3	ug/g	4	58	40 140		
		benzoic acid	<	5.0	ug/g					
		4-chloroaniline		2.6	ug/g	4	64	40 140		
		hexachlorobutadiene		2.5	ug/g	4	62	40 140		
		4-chloro-3-methylphenol		2.7	ug/g	4	68	30 130		
		2-methylnaphthalene		2.27	ug/g	4	57	40 140		
		hexachlorocyclopentadiene	<	1.0	ug/g	4	14	* 40 140		
		2,4,6-trichlorophenol		2.8	ug/g	4	69	30 130		
		2,4,5-trichlorophenol		2.5	ug/g	4	62	30 130		
		2-chloronaphthalene		2.5	ug/g	4	62	40 140		
		2-nitroaniline		2.8	ug/g	4	69	40 140		
		acenaphthylene		2.4	ug/g	4	61	40 140		
		dimethylphthalate		2.7	ug/g	4	68	40 140		
		2,6-dinitrotoluene		2.8	ug/g	4	71	40 140		
		2,4-dinitrotoluene		2.8	ug/g	4	70	40 140		
		acenaphthene		2.5	ug/g	4	61	40 140		
		3-nitroaniline		2.9	ug/g	4	72	40 140		
		2,4-dinitrophenol	<	5.0	ug/g					
		dibenzofuran		2.5	ug/g	4	64	40 140		
		4-nitrophenol		2.0	ug/g	4	50	30 130		
		fluorene		2.8	ug/g	4	69	40 140		
		diethyl phthalate		2.8	ug/g	4	70	40 140		
		4-chlorophenyl phenyl ether		2.8	ug/g	4	69	40 140		
		4-nitroaniline		2.8	ug/g	4	71	40 140		
		4,6-dinitro-2-methylphenol	<	2.0	ug/g					
		azobenzene		2.6	ug/g	4	66	40 140		
		N-nitrosodiphenylamine		3.1	ug/g	4	79	40 140		
		4-bromophenyl phenyl ether		2.7	ug/g	4	67	40 140		
		hexachlorobenzene		2.7	ug/g	4	68	40 140		
		pentachlorophenol		1.9	ug/g	4	48	30 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCS5556	phenanthrene		2.7	ug/g	4	67	40 140		
		anthracene		2.6	ug/g	4	65	40 140		
		carbazole		2.7	ug/g	4	68	40 140		
		di-n-butylphthalate		2.7	ug/g	4	68	40 140		
		fluoranthene		2.7	ug/g	4	67	40 140		
		benzidine	<	3.0	ug/g					
		pyrene		3.6	ug/g	4	89	40 140		
		butyl benzyl phthalate		2.9	ug/g	4	73	40 140		
		benzo(a)anthracene		2.8	ug/g	4	71	40 140		
		chrysene		2.7	ug/g	4	67	40 140		
		3,3'-dichlorobenzidine	<	3.0	ug/g					
		bis(2-ethylhexyl)phthalate		3.1	ug/g	4	79	40 140		
		di-n-octyl phthalate		3.4	ug/g	4	86	40 140		
		benzo(b)fluoranthene		2.8	ug/g	4	69	40 140		
		benzo(k)fluoranthene		3.1	ug/g	4	76	40 140		
		benzo(a)pyrene		2.8	ug/g	4	71	40 140		
		indeno(1,2,3-cd)pyrene		3.1	ug/g	4	77	40 140		
		dibenzo(a,h)anthracene		3.1	ug/g	4	79	40 140		
		benzo(g,h,i)perylene		3.2	ug/g	4	80	40 140		
		2-fluorophenol SUR		59	%			21 100		
		phenol-D5 SUR		54	%			10 102		
		2,4,6-tribromophenol SUR		74	%			10 123		
		nitrobenzene-D5 SUR		58	%			35 114		
		2-fluorobiphenyl SUR		73	%			43 116		
		p-terphenyl-D14 SUR		94	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCSD5556	N-nitrosodimethylamine		2.1	ug/g	4	52	40 140	8	30
		aniline		2.0	ug/g	4	50	40 140	8	30
		phenol		2.3	ug/g	4	58	30 130	1	30
		2-chlorophenol		2.6	ug/g	4	65	30 130	4	30
		bis(2-chloroethyl)ether		2.9	ug/g	4	71	40 140	1	30
		1,3-dichlorobenzene		2.3	ug/g	4	57	40 140	4	30
		1,4-dichlorobenzene		2.4	ug/g	4	60	40 140	4	30
		1,2-dichlorobenzene		2.5	ug/g	4	61	40 140	3	30
		benzyl alcohol		2.3	ug/g	4	57	30 130	5	30
		2-methylphenol		2.7	ug/g	4	67	30 130	8	30
		bis(2-chloroisopropyl) ether		2.4	ug/g	4	59	40 140	3	30
		hexachloroethane		2.3	ug/g	4	58	40 140	3	30
		N-nitroso-di-N-propylamine		2.3	ug/g	4	57	40 140	5	30
		4-methylphenol		2.8	ug/g	4	70	30 130	9	30
		nitrobenzene		2.3	ug/g	4	58	40 140	4	30
		isophorone		2.3	ug/g	4	59	40 140	5	30
		2-nitrophenol		2.2	ug/g	4	56	30 130	4	30
		2,4-dimethylphenol		2.7	ug/g	4	66	30 130	1	30
		bis(2-chloroethoxy)methane		2.5	ug/g	4	63	40 140	6	30
		2,4-dichlorophenol		2.7	ug/g	4	68	30 130	2	30
		1,2,4-trichlorobenzene		2.4	ug/g	4	60	40 140	6	30
		naphthalene		2.4	ug/g	4	60	40 140	4	30
		benzoic acid	<	5.0	ug/g					
		4-chloroaniline		2.7	ug/g	4	68	40 140	6	30
		hexachlorobutadiene		2.6	ug/g	4	65	40 140	5	30
		4-chloro-3-methylphenol		2.8	ug/g	4	69	30 130	2	30
		2-methylnaphthalene		2.35	ug/g	4	59	40 140	3	30
		hexachlorocyclopentadiene	<	1.0	ug/g	4	16	* 40 140	14	30
		2,4,6-trichlorophenol		2.8	ug/g	4	71	30 130	2	30
		2,4,5-trichlorophenol		2.6	ug/g	4	66	30 130	5	30
		2-chloronaphthalene		2.5	ug/g	4	63	40 140	1	30
		2-nitroaniline		2.8	ug/g	4	70	40 140	2	30
		acenaphthylene		2.5	ug/g	4	63	40 140	3	30
		dimethylphthalate		2.8	ug/g	4	70	40 140	3	30
		2,6-dinitrotoluene		2.9	ug/g	4	72	40 140	2	30
		2,4-dinitrotoluene		2.9	ug/g	4	72	40 140	3	30
		acenaphthene		2.5	ug/g	4	62	40 140	1	30
		3-nitroaniline		3.0	ug/g	4	76	40 140	5	30
		2,4-dinitrophenol	<	5.0	ug/g					
		dibenzofuran		2.6	ug/g	4	64	40 140	0	30
		4-nitrophenol		2.1	ug/g	4	53	30 130	7	30
		fluorene		2.7	ug/g	4	69	40 140	0	30
		diethyl phthalate		2.8	ug/g	4	71	40 140	1	30
		4-chlorophenyl phenyl ether		2.8	ug/g	4	69	40 140	1	30
		4-nitroaniline		2.9	ug/g	4	72	40 140	2	30
		4,6-dinitro-2-methylphenol	<	2.0	ug/g					
		azobenzene		2.6	ug/g	4	66	40 140	1	30
		N-nitrosodiphenylamine		3.2	ug/g	4	79	40 140	1	30
		4-bromophenyl phenyl ether		2.7	ug/g	4	68	40 140	1	30
		hexachlorobenzene		2.8	ug/g	4	69	40 140	2	30
		pentachlorophenol		1.9	ug/g	4	48	30 130	0	30

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3546/8270D	LCSD5556	phenanthrene		2.7	ug/g	4	67	40 140	1	30
		anthracene		2.6	ug/g	4	65	40 140	1	30
		carbazole		2.8	ug/g	4	69	40 140	2	30
		di-n-butylphthalate		2.8	ug/g	4	70	40 140	3	30
		fluoranthene		2.7	ug/g	4	68	40 140	2	30
		benzidine	<	3.0	ug/g					
		pyrene		3.8	ug/g	4	94	40 140	5	30
		butyl benzyl phthalate		3.0	ug/g	4	75	40 140	4	30
		benzo(a)anthracene		2.9	ug/g	4	73	40 140	3	30
		chrysene		2.7	ug/g	4	68	40 140	2	30
		3,3'-dichlorobenzidine	<	3.0	ug/g					
		bis(2-ethylhexyl)phthalate		3.3	ug/g	4	82	40 140	4	30
		di-n-octyl phthalate		3.6	ug/g	4	90	40 140	5	30
		benzo(b)fluoranthene		3.0	ug/g	4	74	40 140	7	30
		benzo(k)fluoranthene		3.0	ug/g	4	74	40 140	3	30
		benzo(a)pyrene		2.9	ug/g	4	73	40 140	2	30
		indeno(1,2,3-cd)pyrene		3.1	ug/g	4	77	40 140	0	30
		dibenzo(a,h)anthracene		3.2	ug/g	4	79	40 140	0	30
		benzo(g,h,i)perylene		3.3	ug/g	4	81	40 140	1	30
		2-fluorophenol SUR		59	%			21 100		
		phenol-D5 SUR		53	%			10 102		
		2,4,6-tribromophenol SUR		74	%			10 123		
		nitrobenzene-D5 SUR		58	%			35 114		
		2-fluorobiphenyl SUR		71	%			43 116		
		p-terphenyl-D14 SUR		96	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3550B8270D	BLK5555	naphthalene		<	0.50	ug/g				
		2-methylnaphthalene		<	0.50	ug/g				
		acenaphthylene		<	0.50	ug/g				
		acenaphthene		<	0.50	ug/g				
		dibenzofuran		<	0.50	ug/g				
		fluorene		<	0.50	ug/g				
		phenanthrene		<	0.50	ug/g				
		anthracene		<	0.50	ug/g				
		fluoranthene		<	0.50	ug/g				
		pyrene		<	0.50	ug/g				
		benzo(a)anthracene		<	0.50	ug/g				
		chrysene		<	0.50	ug/g				
		benzo(b)fluoranthene		<	0.50	ug/g				
		benzo(k)fluoranthene		<	0.50	ug/g				
		benzo(a)pyrene		<	0.50	ug/g				
		indeno(1,2,3-cd)pyrene		<	0.50	ug/g				
		dibenzo(a,h)anthracene		<	0.50	ug/g				
		benzo(g,h,i)perylene		<	0.50	ug/g				
		2-fluorobiphenyl SUR		81	%			43	116	
		o-terphenyl SUR		83	%			33	141	
SW3550B8270D	LCS5555	naphthalene		3.3	ug/g	4	83	40	140	
		2-methylnaphthalene		3.3	ug/g	4	81	40	140	
		acenaphthylene		3.2	ug/g	4	81	40	140	
		acenaphthene		3.3	ug/g	4	82	40	140	
		dibenzofuran		<	0.50	ug/g				
		fluorene		3.4	ug/g	4	86	40	140	
		phenanthrene		3.2	ug/g	4	80	40	140	
		anthracene		3.2	ug/g	4	79	40	140	
		fluoranthene		3.1	ug/g	4	78	40	140	
		pyrene		3.4	ug/g	4	86	40	140	
		benzo(a)anthracene		3.3	ug/g	4	82	40	140	
		chrysene		3.0	ug/g	4	76	40	140	
		benzo(b)fluoranthene		4.5	ug/g	4	112	40	140	
		benzo(k)fluoranthene		3.2	ug/g	4	81	40	140	
		benzo(a)pyrene		3.5	ug/g	4	87	40	140	
		indeno(1,2,3-cd)pyrene		3.2	ug/g	4	80	40	140	
		dibenzo(a,h)anthracene		3.3	ug/g	4	82	40	140	
		benzo(g,h,i)perylene		3.3	ug/g	4	82	40	140	
		2-fluorobiphenyl SUR		85	%			43	116	
		o-terphenyl SUR		86	%			33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3550B8270D	MS5555	naphthalene	25123-009	3.4	ug/g	3.90	87	40 140		
		2-methylnaphthalene	25123-009	3.3	ug/g	3.90	84	40 140		
		acenaphthylene	25123-009	3.3	ug/g	3.90	85	40 140		
		acenaphthene	25123-009	3.4	ug/g	3.90	87	40 140		
		dibenzofuran	25123-009	<	0.49	ug/g				
		fluorene	25123-009	3.4	ug/g	3.90	88	40 140		
		phenanthrene	25123-009	3.2	ug/g	3.90	82	40 140		
		anthracene	25123-009	3.1	ug/g	3.90	80	40 140		
		fluoranthene	25123-009	3.4	ug/g	3.90	86	40 140		
		pyrene	25123-009	3.3	ug/g	3.90	86	40 140		
		benzo(a)anthracene	25123-009	3.3	ug/g	3.90	85	40 140		
		chrysene	25123-009	3.2	ug/g	3.90	82	40 140		
		benzo(b)fluoranthene	25123-009	5.0	ug/g	3.90	129	40 140		
		benzo(k)fluoranthene	25123-009	3.7	ug/g	3.90	96	40 140		
		benzo(a)pyrene	25123-009	4.0	ug/g	3.90	101	40 140		
		indeno(1,2,3-cd)pyrene	25123-009	3.7	ug/g	3.90	94	40 140		
		dibenzo(a,h)anthracene	25123-009	3.9	ug/g	3.90	99	40 140		
		benzo(g,h,i)perylene	25123-009	4.0	ug/g	3.90	103	40 140		
		2-fluorobiphenyl SUR	25123-009	88	%			43 116		
		o-terphenyl SUR	25123-009	90	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6010C	BLK5553	Silver		<	0.25	ug/g				
		Arsenic		<	0.50	ug/g				
		Barium		<	2.5	ug/g				
		Beryllium		<	0.20	ug/g				
		Cadmium		<	0.20	ug/g				
		Chromium		<	2.5	ug/g				
		Copper		<	2.5	ug/g				
		Nickel		<	2.5	ug/g				
		Lead		<	0.50	ug/g				
		Antimony		<	0.30	ug/g				
		Selenium		<	2.5	ug/g				
		Thallium		<	0.10	ug/g				
		Zinc		<	2.5	ug/g				
SW3051A6010C	CRM5553	Silver		47	ug/g	38	25.1	51.9		
		Arsenic		460	ug/g	400	292	508		
		Barium		30	ug/g	25	0	51.3		
		Beryllium		<	0.42	ug/g				
		Cadmium			18	ug/g	15	8.71	22	
		Chromium			19	ug/g	14	2.45	24.7	
		Copper			860	ug/g	730	592	866	
		Nickel			16	ug/g	17	6.2	27.5	
		Lead			5300	ug/g	5100	3753	6469	
		Antimony			11	ug/g	8.4	0	21.4	
		Selenium			11	ug/g	6.6	0	18.4	
		Thallium			7.2	ug/g	5.9	0	13.6	
		Zinc			2800	ug/g	3000	2447	3575	
SW3051A6010C	CRMD5553	Silver		48	ug/g	38	25.1	51.9		35
SW3051A6010C	CRMD5553	Arsenic		470	ug/g	400	292	508	3	35
		Barium		33	ug/g	25	0	51.3	11	35
		Beryllium		<	0.42	ug/g				35
		Cadmium			19	ug/g	15	8.71	22	5
		Chromium			19	ug/g	14	2.45	24.7	0
		Copper			850	ug/g	730	592	866	1
		Nickel			16	ug/g	17	6.2	27.5	3
		Lead			5500	ug/g	5100	3753	6469	4
		Antimony			12	ug/g	8.4	0	21.4	11
		Selenium			11	ug/g	6.6	0	18.4	2
		Thallium			6.6	ug/g	5.9	0	13.6	9
		Zinc			2800	ug/g	3000	2447	3575	1
										35

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6010C DUP5553		Silver	25123-002	<	0.32	ug/g				35
		Arsenic	25123-002		13	ug/g			5	35
		Beryllium	25123-002		0.87	ug/g			14	35
		Cadmium	25123-002	<	0.25	ug/g				35
		Chromium	25123-002		27	ug/g			7	35
		Copper	25123-002		35	ug/g			3	35
		Nickel	25123-002		20	ug/g			10	35
		Lead	25123-002		210	ug/g			3	35
		Antimony	25123-002	<	0.38	ug/g				35
		Selenium	25123-002	<	3.2	ug/g				35
		Thallium	25123-002	<	0.13	ug/g				35
		Zinc	25123-002		83	ug/g			5	35
SW3051A6010C MS5553		Silver	25123-002	15	ug/g	15	93	75	125	
		Arsenic	25123-002	42	ug/g	31	88	75	125	
		Beryllium	25123-002	29	ug/g	31	89	75	125	
		Cadmium	25123-002	27	ug/g	31	86	75	125	
		Chromium	25123-002	58	ug/g	31	104	75	125	
		Copper	25123-002	73	ug/g	31	114	75	125	
		Nickel	25123-002	47	ug/g	31	92	75	125	
		Lead	25123-002	250	ug/g	31.7	152	*	75	125
		Antimony	25123-002	25	ug/g	31	77	75	125	
		Selenium	25123-002	27	ug/g	31	79	75	125	
		Thallium	25123-002	30	ug/g	31	95	75	125	
		Zinc	25123-002	110	ug/g	31.7	109	75	125	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3051A6010C	BLK5571	Silver		<	0.25	ug/g					
		Arsenic		<	0.50	ug/g					
		Beryllium		<	0.20	ug/g					
		Cadmium		<	0.20	ug/g					
		Chromium		<	2.5	ug/g					
		Copper		<	2.5	ug/g					
		Nickel		<	2.5	ug/g					
		Lead		<	0.50	ug/g					
		Antimony		<	0.30	ug/g					
		Selenium		<	2.5	ug/g					
		Thallium		<	0.10	ug/g					
		Zinc		<	2.5	ug/g					
SW3051A6010C	CRM5571	Silver		45	ug/g	38		25.1	51.9		
		Arsenic		440	ug/g	400		292	508		
		Barium		26	ug/g	25		0	51.3		
		Beryllium		<	0.42	ug/g					
		Cadmium			20	ug/g	15	8.71	22		
		Chromium			17	ug/g	14	2.45	24.7		
		Copper			850	ug/g	730	592	866		
		Nickel			13	ug/g	17	6.2	27.5		
		Lead			5500	ug/g	5100	3753	6469		
		Antimony			10	ug/g	8.4	0	21.4		
		Selenium			5.8	ug/g	6.6	0	18.4		
		Thallium			4.2	ug/g	5.9	0	13.6		
		Zinc			2600	ug/g	3000	2447	3575		
SW3051A6010C	CRMD5571	Silver		44	ug/g	38		25.1	51.9	1	35
		Arsenic		480	ug/g	400		292	508	7	35
		Barium		25	ug/g	25		0	51.3	7	35
		Beryllium		<	0.42	ug/g					35
		Cadmium			19	ug/g	15	8.71	22	6	35
		Chromium			18	ug/g	14	2.45	24.7	6	35
		Copper			900	ug/g	730	592	866	5	35
		Nickel			16	ug/g	17	6.2	27.5	23	35
		Lead			4700	ug/g	5100	3753	6469	16	35
		Antimony			8.2	ug/g	8.4	0	21.4	24	35
		Selenium			5.8	ug/g	6.6	0	18.4	0	35
		Thallium			1.6	ug/g	5.9	0	13.6	90	*
		Zinc			2700	ug/g	3000	2447	3575	6	35
SW3051A6010C	DUP5571	Silver	25123-016	<	0.35	ug/g					35
		Arsenic	25123-016		2.5	ug/g			24		35
		Barium	25123-016		18	ug/g			19		35
		Cadmium	25123-016	<	0.28	ug/g					35
		Chromium	25123-016		8.9	ug/g			23		35
		Lead	25123-016		4.3	ug/g			20		35
		Selenium	25123-016	<	3.5	ug/g					35

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3051A6010C	MS5571	Silver	25123-016	15	ug/g	17	86	75 125		
		Arsenic	25123-016	35	ug/g	34	92	75 125		
		Barium	25123-016	56	ug/g	34	98	75 125		
		Cadmium	25123-016	33	ug/g	34	93	75 125		
		Chromium	25123-016	44	ug/g	34	94	75 125		
		Lead	25123-016	35	ug/g	34	86	75 125		
		Selenium	25123-016	31	ug/g	34	85	75 125		
SW7471B	BLK5564	Mercury		<	ug/g					
SW7471B	CRM5564	Mercury		2.5	ug/g	2.2		0.98 3.52		
SW7471B	CRMD5564	Mercury		1.2	ug/g	1.1		0.49 1.76		35
SW7471B	DUP5564	Mercury	25094-001	<	ug/g					35
SW7471B	MS5564	Mercury	25094-001	0.48	ug/g	0.417	110	75 125		
SW7471B	MS5564	Mercury	25123-010	0.76	ug/g	0.382	129	75 125		
SW7471B	BLK5565	Mercury		<	ug/g					
SW7471B	CRM5565	Mercury		1.3	ug/g	1.1		0.49 1.76		
SW7471B	CRMD5565	Mercury		1.5	ug/g	1.1		0.49 1.76	12	35
SW7471B	DUP5565	Mercury	25123-024	2.4	ug/g					35
SW7471B	MS5565	Mercury	25123-024	4.7	ug/g	0.39	1104	75 125		

File : X:\SVOA01\2012\OCT12\100912\S1100938.D

Operator : jlz

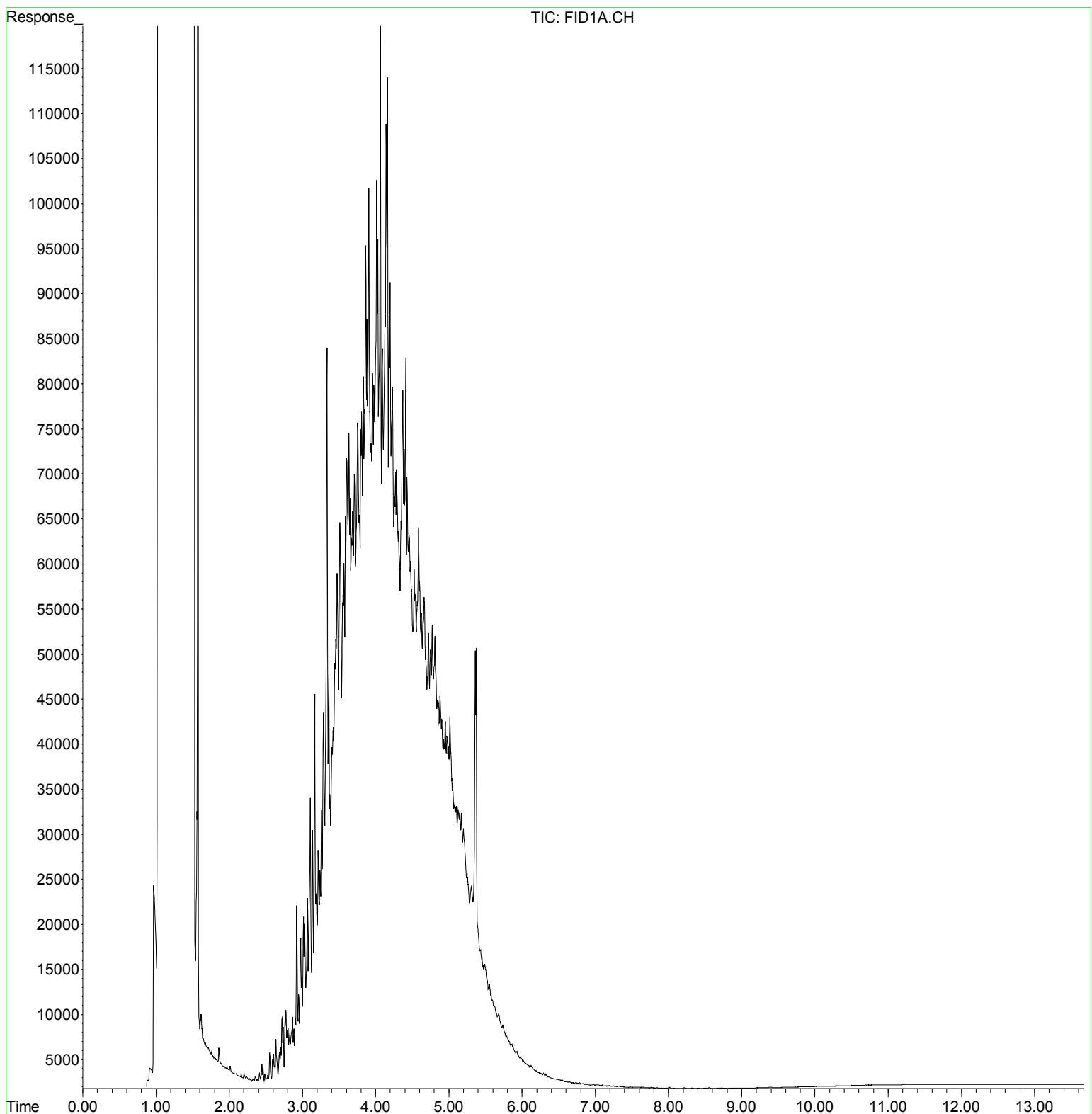
Acquired : 10 Oct 2012 12:06 am using AcqMethod S1ZIPALI.M

Instrument : SVOA01

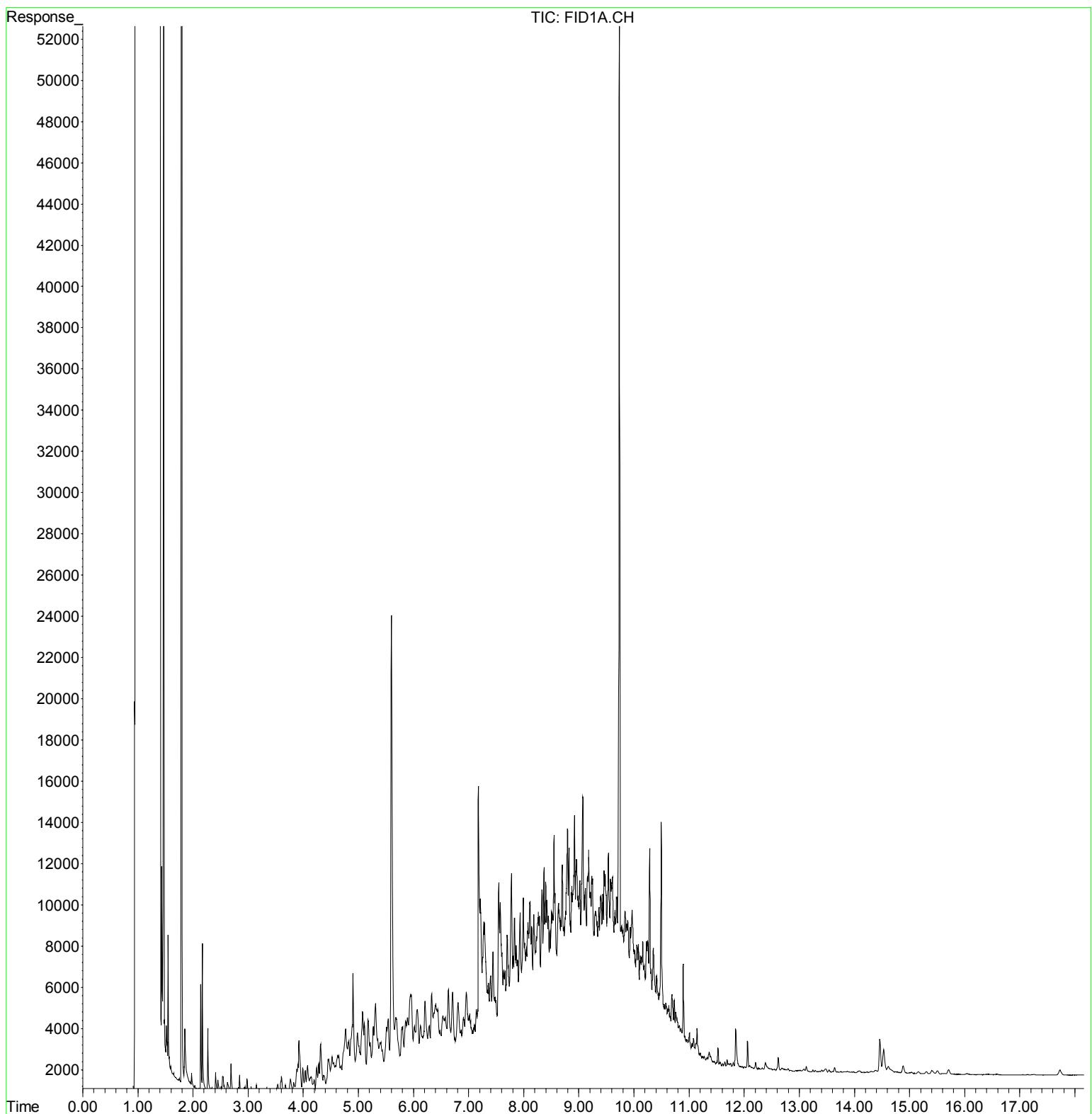
Sample Name: 25123-15 ali x1

Misc Info : Field ID: CA-SB-12 15-18 [Aliphatic portion]

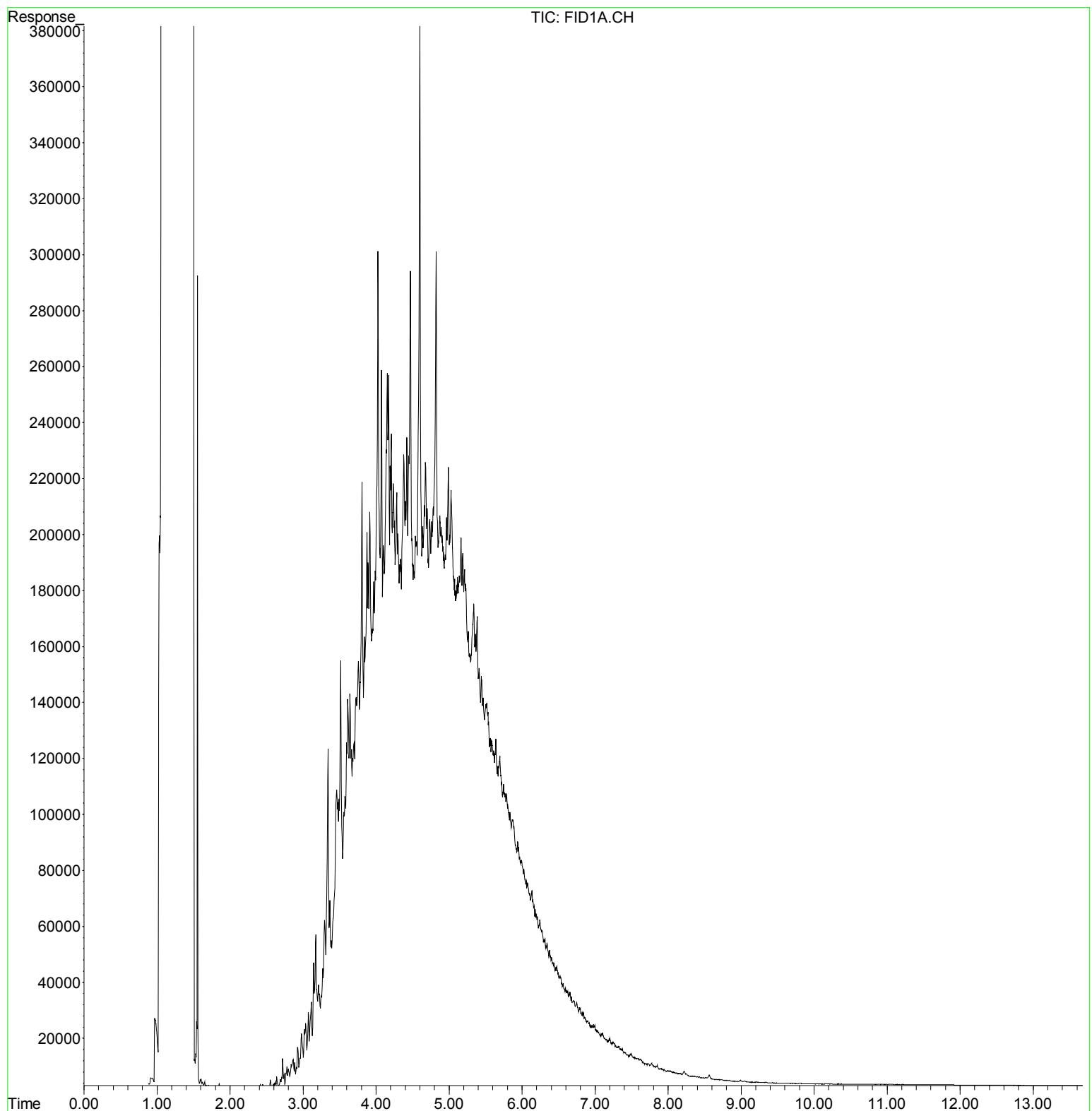
Vial Number: 38



File : X:\SVOA01\2012\OCT12\101412\S1101421.D
Operator : jlz
Acquired : 14 Oct 2012 10:28 pm using AcqMethod S1ZIPARO.M
Instrument : SVOA01
Sample Name: 25123-15 aro x1
Misc Info : Field ID: CA-SB-12 15-18 [Aromatic portion]
Vial Number: 21



File : X:\SVOA01\2012\OCT12\100912\S1100941.D
Operator : jlz
Acquired : 10 Oct 2012 1:00 am using AcqMethod S1ZIPALI.M
Instrument : SVOA01
Sample Name: 25123-16 ali x1
Misc Info : Field ID: CA-SB-13 10-12 [Aliphatic portion]
Vial Number: 41



File : X:\SVOA01\2012\OCT12\100912\S1100942.D

Operator : jlz

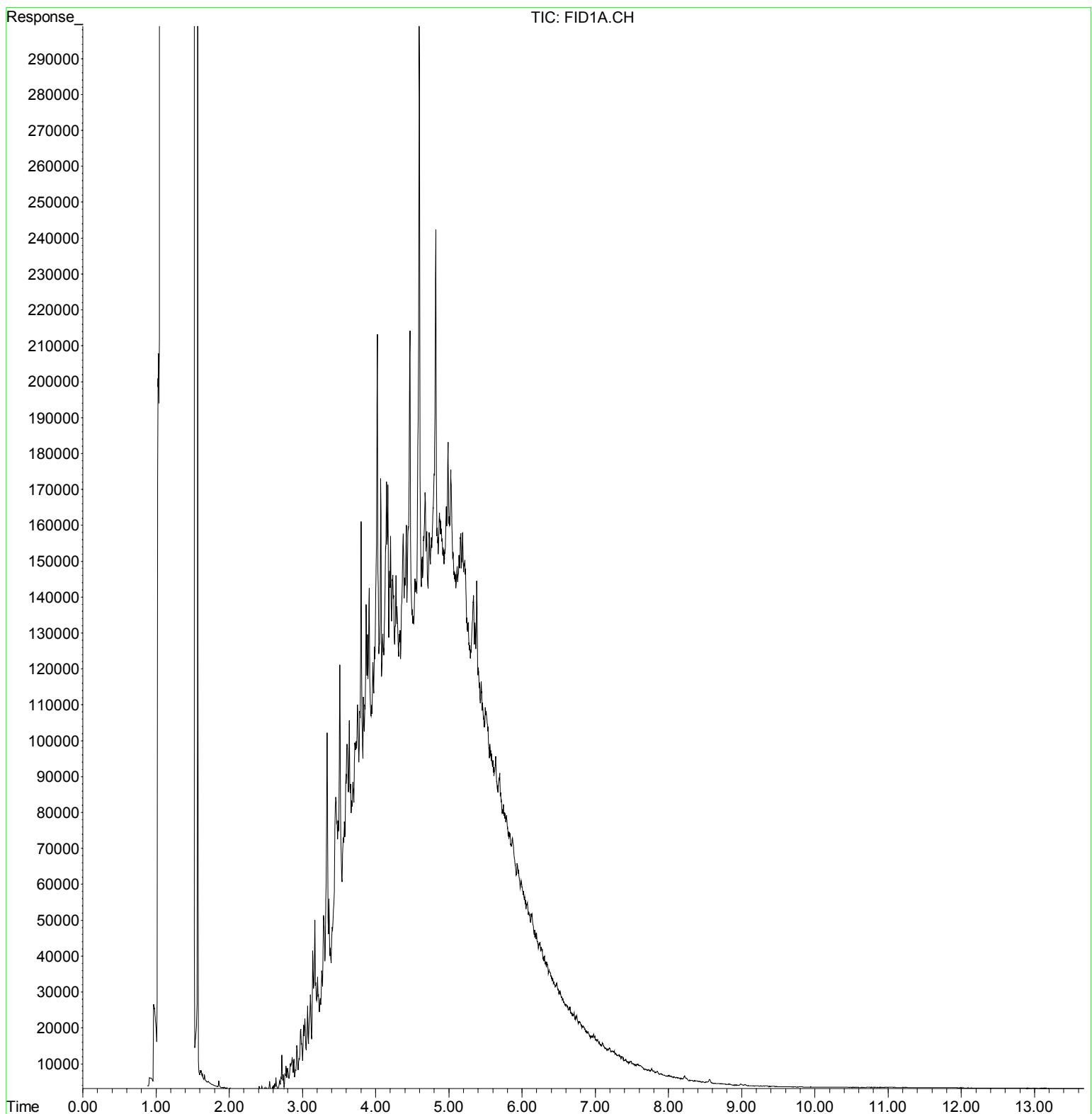
Acquired : 10 Oct 2012 1:19 am using AcqMethod S1ZIPALI.M

Instrument : SVOA01

Sample Name: 25123-17 ali x1

Misc Info : Field ID: SB-DUP-1 [Aliphatic portion]

Vial Number: 42



File : X:\SVOA01\2012\OCT12\101412\S1101423.D

Operator : jlz

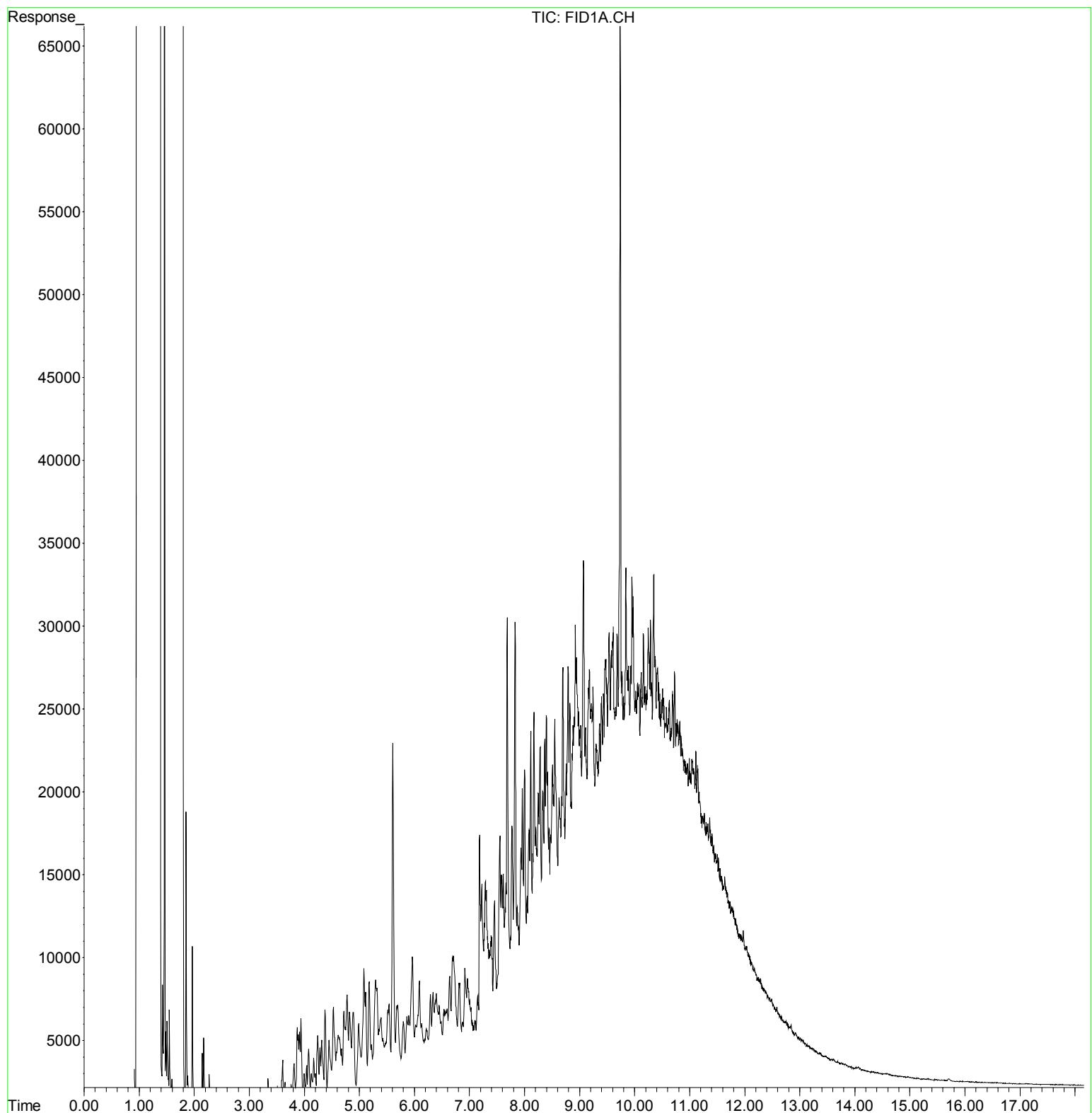
Acquired : 14 Oct 2012 11:13 pm using AcqMethod S1ZIPARO.M

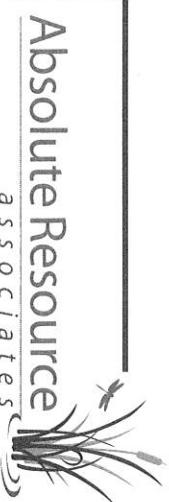
Instrument : SVOA01

Sample Name: 25123-17 aro x1

Misc Info : Field ID: SB-DUP-1 [Aromatic portion]

Vial Number: 23





124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST**

25123

PAGE 1 OF 5

Absolute Resource

a s s o c i a t e s

124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

25123

Absolute Resource

associates

absoluteresourceassociates.com

124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

25123

ANALYSIS REQUEST

Company Name: CREDENCE ASSOCIATES	Project Name: MERC
Company Address: 116 MAIN ST WESTBROOK ME 04092	Project #: 1001150
Report To: JUDD NEWCOMB	Project Location: NH MA VT Other
Phone #: 207-332-5387	Protocol: <input checked="" type="checkbox"/> CEM <input type="checkbox"/> MCP <input type="checkbox"/> SDWA <input type="checkbox"/> NPDES <input type="checkbox"/> OTHER
Invoice To: Same	Reporting QAPP GW-1 S-1
	Limits: EPA DW Other
	Quote #: CREDENCE <input type="checkbox"/> NH GREE/ODD Fund Pricing
	PO #: 1001150

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix	Preservation Method	Sampling		SAMPLER																						
					WATER	SOLID		OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE	TIME													
25123-12	CA-SB-10	10-15	2		X					4/12	10/13/12	1030	TR	X	X	VOC 8260	<input type="checkbox"/>	VOC 8260 NHDES	<input type="checkbox"/>	VOC 8260 MADEP									
-13	CA-SB-11	10-15	3														<input type="checkbox"/>	VOC 624	<input type="checkbox"/>	VOC BTEX	<input type="checkbox"/>	MIBE, only <input type="checkbox"/>	<input type="checkbox"/>	VOC 8021VT					
-14	SB-DUP-2	15	3														X	VPH MADEP	<input type="checkbox"/>	MEGRO	<input type="checkbox"/>	GRO 8015							
-15	CA-SB-12	10-12	1														<input type="checkbox"/>	VOC 524.2	<input type="checkbox"/>	VOC 524.2 NH List	<input type="checkbox"/>	Gases-List:							
-16	CA-SB-13	10-12	1														<input type="checkbox"/>	TPH	<input type="checkbox"/>	DRO 8015	<input type="checkbox"/>	MEDRO	<input checked="" type="checkbox"/>	EPH MADEP	<input type="checkbox"/>	TPH Fingerprint			
-17	SB-DUP-1	1															<input type="checkbox"/>	8270PAH	<input checked="" type="checkbox"/>	8270ABN	<input type="checkbox"/>	625	<input type="checkbox"/>	EDB 504.1					
-18	CA-SB-14	0-3	2														<input type="checkbox"/>	8082 PCB	<input type="checkbox"/>	8081 Pesticides	<input type="checkbox"/>	608 Pest/PCB	3540 EXT						
-19	CA-SB-15	5-5	2														<input type="checkbox"/>	O&G 1664	<input type="checkbox"/>	Mineral O&G SM5520F									
-20	CA-SB-16	5-5	2														<input type="checkbox"/>	pH	<input type="checkbox"/>	BOD	<input type="checkbox"/>	Conductivity	<input type="checkbox"/>	Turbidity					
-21	CA-SB-17	1	1														<input type="checkbox"/>	TSS	<input type="checkbox"/>	TDS	<input type="checkbox"/>	TS	<input type="checkbox"/>	TVS	<input type="checkbox"/>	Alkalinity			
-22	CA-SB-18	1	1														<input type="checkbox"/>	RCRA Metals	<input checked="" type="checkbox"/>	Priority Pollutant Metals	<input type="checkbox"/>	TAL Metals							
																	<input type="checkbox"/>	Total Metals-list:											
																	<input type="checkbox"/>	Dissolved Metals-list:											
																	<input type="checkbox"/>	Ammonia	<input type="checkbox"/>	COD	<input type="checkbox"/>	TKN	<input type="checkbox"/>	TN	<input type="checkbox"/>	TON	<input type="checkbox"/>	TOC	
																	<input type="checkbox"/>	T-Phosphorus	<input type="checkbox"/>	Phenols	<input type="checkbox"/>	Bacteria P/A	<input type="checkbox"/>	Bacteria MPN					
																	<input type="checkbox"/>	Cyanide	<input type="checkbox"/>	Sulfide	<input type="checkbox"/>	Nitrate + Nitrite	<input type="checkbox"/>	Ortho P					
																	<input type="checkbox"/>	Nitrate	<input type="checkbox"/>	Nitrite	<input type="checkbox"/>	Chloride	<input type="checkbox"/>	Sulfate	<input type="checkbox"/>	Bromide	<input type="checkbox"/>	Fluoride	
																	<input type="checkbox"/>	Corrosivity	<input type="checkbox"/>	Reactive CN	<input type="checkbox"/>	Reactive S-	<input type="checkbox"/>	Ignitability/FP					
																	<input type="checkbox"/>	TCLP Metals	<input type="checkbox"/>	TCLP VOC	<input type="checkbox"/>	TCLP SVOC	<input type="checkbox"/>	TCLP Pesticide					
																	<input type="checkbox"/>	Subcontract:	<input type="checkbox"/>	Grain Size	<input type="checkbox"/>	Herbicides	<input type="checkbox"/>	Formaldehyde					
																	ICMA 8 metals												
																	VPH/EPH TARGET NAMES ONLY												
																	PAH 8d70												
																	Grab (G) or Composite (C)												

TAT REQUESTED

Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)
*Date Needed **10/10/12**

SPECIAL INSTRUCTIONS

See absoluteresourceassociates.com
for sample acceptance policy and
current accreditation lists.

REPORTING INSTRUCTIONS

PDF (e-mail address) **INVERNESS@CREDENCELLC.COM**

HARD COPY REQUIRED

FAX (FAX#) **1001150**

OTHER (specify)

Relinquished by Sample#: **10/12/12** Date **10/15** Time **14:00** Received by: **CREDENCE Lab STAFF** Date **10/12/12** Time **16:15**

Relinquished by: **Chad Stoenke** Received by: **Chad Stoenke** Date **10/12/12** Time **11:30** Received by Laboratory: **Chad Stoenke** Date **10/12/12** Time **11:30** Way Bill#: **1001150**

CUSTODY RECORD

Relinquished by: **Chad Stoenke** Received by: **Chad Stoenke** Date **10/12/12** Time **11:30**

OSD-01 Revision 10/6/11

Absolute Resource

associates

124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

25123

PAGE 3 OF 3

Company Name:
CREDENCE ASSOCIATES
Company Address:
776 MAIN ST WESTBROOK ME 04092

Report To:
JUDD NEWCOMB
Phone #:
207-832-5387

Invoice To:
Same

ANALYSIS REQUEST

Project Name:	MERC
Project #:	12001150
Project Location:	NH MA VT Other
Protocol:	<input checked="" type="checkbox"/> PCB <input type="checkbox"/> SDWA <input type="checkbox"/> NPDES <input type="checkbox"/> MCP <input type="checkbox"/> NHDES <input type="checkbox"/> OTHER
Reporting Limits:	QAPP GW-1 S-1 EPA DW Other
Quote #:	CREDENCE <input type="checkbox"/> NH GREE/ODD PO # 12001150 <input type="checkbox"/> Fund Pricing

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix	Preservation Method	Sampling		Grab (G) or Composite (C)
					DATE	TIME	
25-23	CA-SS-4	1	WATER	X	10/3/12	1045 Tr	X
-24	CA-SS-5	1	SOLID		10/3/12	1300 Tr	
-25	CA-CC-1	2	OTHER	X	10/3/12	1300 Tr	X
-26	CA-CC-2	2	HCl	X	10/3/12	1300 Tr	X
-27	CA-CC-3	2	HNO ₃	X	10/3/12	1300 Tr	X
-28	CA-CC-4	1	H ₂ SO ₄	X	10/3/12	1300 Tr	X
-29	CA-CC-5	1	NaOH	X	10/3/12	1300 Tr	X
-30	STACK	1	MeOH	X	10/3/12	1300 Tr	X
-31	TRIP BANK	1	OTHER (Specify)	X	10/3/12	1300 Tr	X

TAT REQUESTED

Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)

*Date Needed **10/10/12**

See absoluteresourceassociates.com
for sample acceptance policy and
current accreditation lists.

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS PDF (e-mail address) **JNewcombs.CredenceLLC.com**

HARD COPY REQUIRED FAX (FAX#) _____

OTHER (specify) _____

RECEIVED ON ICE YES NO
TEMPERATURE **0** °C

Relinquished by:

John M. Judd

Date **10/3/12**

Time **1015**

Received by: **CREDENCE COLD STORAGE**

Received by: *John M. Judd*

Date **10/11/12**

Time **1130**

Received by Laboratory: *John M. Judd*

Received by Laboratory: *John M. Judd*

Date **10/11/12**

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Received by Laboratory: *John M. Judd*

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Date **10/11/12**

Time **1130**

Received by Laboratory: *John M. Judd*

Date **1**

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

Judd Newcomb
CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: None
Job ID: 25149
Date Received: 10/10/12

Project: MERC 12001151

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

Cliff Chase
Partner, Technical Director

Date of Approval: 10/23/2012
Total number of pages: 62

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-MW-3	Water	10/9/2012 15:20	25149-001	EPH in water by MADEP Method Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Barium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Mercury in water by 7470 Lead in water by 6010 Selenium in water by 6010
CA-MW-6	Water	10/9/2012 13:20	25149-002	Acid & Base/Neutral Extractables in water by 8270 EPH in water by MADEP Method Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Beryllium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Antimony in water by 6010 Selenium in water by 6010 Thallium in water by 6010 Zinc in water by 6010 VOCs in water by 8260 VPH in water by MA DEP Method
CA-MW-7	Water	10/9/2012 14:15	25149-003	Acid & Base/Neutral Extractables in water by 8270 EPH in water by MADEP Method Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Beryllium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Antimony in water by 6010 Selenium in water by 6010 Thallium in water by 6010 Zinc in water by 6010 VOCs in water by 8260 VPH in water by MA DEP Method
CA-MW-8	Water	10/9/2012 12:12	25149-004	PAHs in water by 8270 Water Digestion for ICP Analysis

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-MW-8	Water	10/9/2012 12:12	25149-004	Silver in water by 6010 Arsenic in water by 6010 Beryllium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Antimony in water by 6010 Selenium in water by 6010 Thallium in water by 6010 Zinc in water by 6010
CA-MW-9	Water	10/9/2012 10:53	25149-005	Acid & Base/Neutral Extractables in water by 8270 Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Beryllium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Antimony in water by 6010 Selenium in water by 6010 Thallium in water by 6010 Zinc in water by 6010 VOCs in water by 8260
CA-MW-10	Water	10/9/2012 9:42	25149-006	Acid & Base/Neutral Extractables in water by 8270 Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Barium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010 Mercury in water by 7470 Lead in water by 6010 Selenium in water by 6010 VOCs in water by 8260
CA-MW-DUP-1	Water	10/9/2012 0:00	25149-007	Acid & Base/Neutral Extractables in water by 8270 EPH in water by MADEP Method Water Digestion for ICP Analysis Silver in water by 6010 Arsenic in water by 6010 Beryllium in water by 6010 Cadmium in water by 6010 Chromium in water by 6010

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-MW-DUP-1	Water	10/9/2012 0:00	25149-007	Copper in water by 6010 Mercury in water by 7470 Nickel in water by 6010 Lead in water by 6010 Antimony in water by 6010 Selenium in water by 6010 Thallium in water by 6010 Zinc in water by 6010 VOCs in water by 8260 VPH in water by MA DEP Method
Trip Blank	Water	10/9/2012 0:00	25149-008	VOCs in water by 8260
Trip Blank	Water	10/9/2012 0:00	25149-009	VPH in water by MA DEP Method

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
vinyl chloride	2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
cis-1,2-dichloroethene	2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
Surrogate Recovery										
dibromofluoromethane SUR	100	78-114	%	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
toluene-D8 SUR	101	88-110	%	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
4-bromofluorobenzene SUR	100	86-115	%	1	LMM	1202467	10/15/12	17:57	SW5030B8260B	
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2,4-trimethylbenzene	2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
4-isopropyltoluene	14	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
Surrogate Recovery										
dibromofluoromethane SUR	103	78-114	%	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
toluene-D8 SUR	103	88-110	%	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
4-bromofluorobenzene SUR	100	86-115	%	1	LMM	1202467	10/15/12	18:29	SW5030B8260B	
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-005

Sample ID: CA-MW-9

Matrix: Water

Sampled: 10/9/12 10:53

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-005

Sample ID: CA-MW-9

Matrix: Water

Sampled: 10/9/12 10:53

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
Surrogate Recovery										
dibromofluoromethane SUR	103	78-114	%	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
toluene-D8 SUR	102	88-110	%	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
4-bromofluorobenzene SUR	102	86-115	%	1	LMM	1202467	10/15/12	19:00	SW5030B8260B	
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-006

Sample ID: CA-MW-10

Matrix: Water

Sampled: 10/9/12 9:42

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
benzene	23	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
chlorobenzene	270	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-006

Sample ID: CA-MW-10

Matrix: Water

Sampled: 10/9/12 9:42

Parameter	Result	Reporting Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Analysis		
							Batch	Date	Time
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
Surrogate Recovery						Limits			
dibromofluoromethane SUR	99	78-114	%	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
toluene-D8 SUR	99	88-110	%	1	LMM	1202467	10/15/12	19:32	SW5030B8260B
4-bromofluorobenzene SUR	97	86-115	%	1	LMM	1202467	10/15/12	19:32	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
cis-1,2-dichloroethene	2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
Surrogate Recovery										
dibromofluoromethane SUR	99	78-114	%	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
toluene-D8 SUR	102	88-110	%	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	
4-bromofluorobenzene SUR	98	86-115	%	1	LMM	1202467	10/15/12	20:03	SW5030B8260B	

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-008

Sample ID: Trip Blank

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Result	Reporting Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Analysis		
							Batch	Date	Time
dichlorodifluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
chloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
vinyl chloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
bromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
chloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
trichlorofluoromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
diethyl ether	< 5	5	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
acetone	< 50	50	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1-dichloroethene	< 1	1	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
methylene chloride	< 5	5	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
carbon disulfide	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
trans-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
2-butanone (MEK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
2,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
cis-1,2-dichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
chloroform	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
bromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
tetrahydrofuran (THF)	< 10	10	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1,1-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
carbon tetrachloride	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2-dichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
benzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
trichloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
bromodichloromethane	< 1	1	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
dibromomethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
4-methyl-2-pentanone (MIBK)	< 10	10	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
cis-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
toluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
trans-1,3-dichloropropene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
2-hexanone	< 10	10	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1,2-trichloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,3-dichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
tetrachloroethene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
dibromochloromethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2-dibromoethane (EDB)	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
chlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1,1,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
ethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
m&p xylenes	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-008

Sample ID: Trip Blank

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis		
	Result	Limit					Batch	Date	Time
o-xylene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
styrene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
bromoform	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
isopropylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,1,2,2-tetrachloroethane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2,3-trichloropropane	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
n-propylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
bromobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,3,5-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
2-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
4-chlorotoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
tert-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2,4-trimethylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
sec-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,3-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
4-isopropyltoluene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,4-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2-dichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
n-butylbenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2-dibromo-3-chloropropane (DBCP)	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2,4-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
hexachlorobutadiene	< 0.5	0.5	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
naphthalene	< 5	5	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
1,2,3-trichlorobenzene	< 2	2	ug/L	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
Surrogate Recovery									
dibromofluoromethane SUR	101	78-114	%	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
toluene-D8 SUR	100	88-110	%	1	LMM	1202467	10/15/12	13:43	SW5030B8260B
4-bromofluorobenzene SUR	101	86-115	%	1	LMM	1202467	10/15/12	13:43	SW5030B8260B

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Received on ice at 4°C, in satisfactory condition.

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
benzene	< 1	1	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
toluene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
ethylbenzene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
m&p-xylenes	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
o-xylene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
naphthalene	< 5	5	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
C9-C10 Aromatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:06	MA VPH	
Surrogate Recovery										
2,5-dibromotoluene as Aromatic SUR		116	70-130	%	1	LMM	1202475	10/15/12	18:06	MA VPH
2,5-dibromotoluene as Aliphatic SUR		115	70-130	%	1	LMM	1202475	10/15/12	18:06	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Received on ice at 4°C, in satisfactory condition.

Sampled: 10/9/12 14:15

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
Unadjusted C9-C12 Aliphatics	110	100	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
benzene	< 1	1	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
toluene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
ethylbenzene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
m&p-xylenes	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
o-xylene	< 2	2	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
naphthalene	< 5	5	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
C9-C10 Aromatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	18:39	MA VPH	
Surrogate Recovery		Limits								
2,5-dibromotoluene as Aromatic SUR	111	70-130	%	1	LMM	1202475	10/15/12	18:39	MA VPH	
2,5-dibromotoluene as Aliphatic SUR	99	70-130	%	1	LMM	1202475	10/15/12	18:39	MA VPH	

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Received on ice at 4°C, in satisfactory condition.

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
benzene	< 1	1	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
toluene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
ethylbenzene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
m&p-xylenes	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
o-xylene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
naphthalene	< 5	5	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
C9-C10 Aromatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:12	MA VPH	
Surrogate Recovery										
2,5-dibromotoluene as Aromatic SUR		112	70-130	%	1	LMM	1202475	10/15/12	19:12	MA VPH
2,5-dibromotoluene as Aliphatic SUR		112	70-130	%	1	LMM	1202475	10/15/12	19:12	MA VPH

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-009

Sample ID: Trip Blank

Matrix: Water

Received on ice at 4°C, in satisfactory condition.

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr Dil'n		Analyst	Prep Date	Analysis			Reference
	Result	Limit	Units	Factor			Batch	Date	Time	
Unadjusted C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
Unadjusted C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
methyl t-butyl ether (MTBE)	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
benzene	< 1	1	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
toluene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
ethylbenzene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
m&p-xylenes	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
o-xylene	< 2	2	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
naphthalene	< 5	5	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
C5-C8 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
C9-C12 Aliphatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
C9-C10 Aromatics	< 100	100	ug/L	1	LMM	1202475	10/15/12	19:45	MA VPH	
Surrogate Recovery		Limits								
2,5-dibromotoluene as Aromatic SUR	110	70-130	%	1	LMM	1202475	10/15/12	19:45	MA VPH	
2,5-dibromotoluene as Aliphatic SUR	110	70-130	%	1	LMM	1202475	10/15/12	19:45	MA VPH	

Hydrocarbon Range data exclude concentrations of any surrogate(s) and/or internal standards eluting in that range.

C5-C8 Aliphatic Hydrocarbons exclude the concentration of target analytes eluting in that range.

C9-C12 Aliphatic Hydrocarbons exclude concentration of target analytes eluting in that range AND C9-C10 Aromatics.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
N-nitrosodimethylamine	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
aniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
phenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
bis(2-chloroethyl)ether	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
1,3-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
1,4-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
1,2-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzyl alcohol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
bis(2-chloroisopropyl) ether	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
hexachloroethane	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
N-nitroso-di-N-propylamine	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
nitrobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
isophorone	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
bis(2-chloroethoxy)methane	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
1,2,4-trichlorobenzene	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
naphthalene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
4-chloroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
hexachlorobutadiene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
hexachlorocyclopentadiene	< 10	10	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
2-chloronaphthalene	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2-nitroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
acenaphthylene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
dimethylphthalate	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2,6-dinitrotoluene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2,4-dinitrotoluene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
acenaphthene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
3-nitroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
dibenzofuran	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD 10/12/12	5563	10/16/12	20:14	SW3510C8270D
fluorene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D
diethyl phthalate	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	20:51	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
4-chlorophenyl phenyl ether	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
4-nitroaniline	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	10/12/12	5563	10/16/12	20:14	SW3510C8270D
azobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
N-nitrosodiphenylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
4-bromophenyl phenyl ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
hexachlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
pentachlorophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	20:14	SW3510C8270D
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
carbazole	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
di-n-butylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
butyl benzyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
3,3'-dichlorobenzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
bis(2-ethylhexyl)phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
di-n-octyl phthalate	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
dibenz(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
Surrogate Recovery										
2-fluorophenol SUR	27	21-100	%	1	AJD	10/12/12	5563	10/16/12	20:14	SW3510C8270D
phenol-D5 SUR	16	10-102	%	1	AJD	10/12/12	5563	10/16/12	20:14	SW3510C8270D
2,4,6-tribromophenol SUR	87	10-123	%	1	AJD	10/12/12	5563	10/16/12	20:14	SW3510C8270D
nitrobenzene-D5 SUR	69	35-114	%	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
2-fluorobiphenyl SUR	79	43-116	%	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D
p-terphenyl-D14 SUR	77	33-141	%	1	AJD	10/12/12	5563	10/16/12	20:51	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			
	Result	Limit				Batch	Date	Time	Reference
N-nitrosodimethylamine	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
aniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
phenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
bis(2-chloroethyl)ether	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
1,3-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
1,4-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
1,2-dichlorobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzyl alcohol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
bis(2-chloroisopropyl) ether	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
hexachloroethane	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
N-nitroso-di-N-propylamine	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
nitrobenzene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
isophorone	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
bis(2-chloroethoxy)methane	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
1,2,4-trichlorobenzene	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
naphthalene	1.6	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
4-chloroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
hexachlorobutadiene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
2-methylnaphthalene	3.3	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
hexachlorocyclopentadiene	< 10	10	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
2-chloronaphthalene	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2-nitroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
acenaphthylene	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
dimethylphthalate	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2,6-dinitrotoluene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2,4-dinitrotoluene	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
acenaphthene	0.9	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
3-nitroaniline	< 2	2	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
dibenzofuran	< 0.5	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD 10/12/12	5563	10/16/12	21:29	SW3510C8270D
fluorene	0.8	0.5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D
diethyl phthalate	< 5	5	ug/L	1	AJD 10/12/12	5563	10/16/12	22:06	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
4-chlorophenyl phenyl ether	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
4-nitroaniline	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	10/12/12	5563	10/16/12	21:29	SW3510C8270D
azobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
N-nitrosodiphenylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
4-bromophenyl phenyl ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
hexachlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
pentachlorophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	21:29	SW3510C8270D
phenanthrene	0.8	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
carbazole	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
di-n-butylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
butyl benzyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
3,3'-dichlorobenzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
bis(2-ethylhexyl)phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
di-n-octyl phthalate	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
dibenz(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
Surrogate Recovery										
2-fluorophenol SUR	31	21-100	%	1	AJD	10/12/12	5563	10/16/12	21:29	SW3510C8270D
phenol-D5 SUR	19	10-102	%	1	AJD	10/12/12	5563	10/16/12	21:29	SW3510C8270D
2,4,6-tribromophenol SUR	112	10-123	%	1	AJD	10/12/12	5563	10/16/12	21:29	SW3510C8270D
nitrobenzene-D5 SUR	70	35-114	%	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
2-fluorobiphenyl SUR	87	43-116	%	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D
p-terphenyl-D14 SUR	80	33-141	%	1	AJD	10/12/12	5563	10/16/12	22:06	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-004

Sample ID: CA-MW-8

Matrix: Water

Sampled: 10/9/12 12:12

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
naphthalene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
2-methylnaphthalene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
acenaphthylene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
acenaphthene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
dibenzofuran	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
fluorene	5.7	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
phenanthrene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
anthracene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
fluoranthene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
pyrene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
benzo(a)anthracene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
chrysene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
benzo(b)fluoranthene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
benzo(k)fluoranthene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
benzo(a)pyrene	< 1.0	1.0	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
indeno(1,2,3-cd)pyrene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
dibenzo(a,h)anthracene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
benzo(g,h,i)perylene	< 2.6	2.6	ug/L	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
Surrogate Recovery										
2-fluorobiphenyl SUR	86	43-116	%	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D
o-terphenyl SUR	87	33-141	%	5	AJD	10/12/12	5563	10/18/12	11:40	SW3510C8270D

Note: Dilution was required due to interferences from hydrocarbons.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-005

Sample ID: CA-MW-9

Matrix: Water

Sampled: 10/9/12 10:53

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Batch	Analysis		
								Date	Time	Reference
N-nitrosodimethylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
aniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
phenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
bis(2-chloroethyl)ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
1,3-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
1,4-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
1,2-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzyl alcohol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
bis(2-chloroisopropyl) ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
hexachloroethane	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
N-nitroso-di-N-propylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
nitrobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
isophorone	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
bis(2-chloroethoxy)methane	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
1,2,4-trichlorobenzene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
4-chloroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
hexachlorobutadiene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
hexachlorocyclopentadiene	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2-chloronaphthalene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
dimethylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2,6-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2,4-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
3-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
dibenzofuran	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
diethyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-005

Sample ID: CA-MW-9

Matrix: Water

Sampled: 10/9/12 10:53

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
4-chlorophenyl phenyl ether	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4-nitroaniline	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
azobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
N-nitrosodiphenylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
4-bromophenyl phenyl ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
hexachlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
pentachlorophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
carbazole	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
di-n-butylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
butyl benzyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
3,3'-dichlorobenzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
bis(2-ethylhexyl)phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
di-n-octyl phthalate	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
dibenzo(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
Surrogate Recovery										
2-fluorophenol SUR	30	21-100	%	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
phenol-D5 SUR	17	10-102	%	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
2,4,6-tribromophenol SUR	99	10-123	%	1	AJD	10/12/12	5563	10/16/12	22:43	SW3510C8270D
nitrobenzene-D5 SUR	76	35-114	%	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
2-fluorobiphenyl SUR	95	43-116	%	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D
p-terphenyl-D14 SUR	98	33-141	%	1	AJD	10/12/12	5563	10/16/12	23:21	SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-006

Sample ID: CA-MW-10

Matrix: Water

Sampled: 10/9/12 9:42

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Batch	Analysis			Reference
								Date	Time		
N-nitrosodimethylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
aniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
phenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
bis(2-chloroethyl)ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
1,3-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
1,4-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
1,2-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
benzyl alcohol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
bis(2-chloroisopropyl) ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
hexachloroethane	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
N-nitroso-di-N-propylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
nitrobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
isophorone	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
bis(2-chloroethoxy)methane	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
1,2,4-trichlorobenzene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
4-chloroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
hexachlorobutadiene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
hexachlorocyclopentadiene	< 10	10	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
2-chloronaphthalene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
dimethylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2,6-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2,4-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
3-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
dibenzofuran	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58		SW3510C8270D
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D
diethyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35		SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-006

Sample ID: CA-MW-10

Matrix: Water

Sampled: 10/9/12 9:42

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Batch	Analysis			Reference
								Date	Time		
4-chlorophenyl phenyl ether	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
4-nitroaniline	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58	SW3510C8270D	
azobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
N-nitrosodiphenylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
4-bromophenyl phenyl ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
hexachlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
pentachlorophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/16/12	23:58	SW3510C8270D	
phenanthrene	0.6	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
carbazole	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
di-n-butylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
butyl benzyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
3,3'-dichlorobenzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
bis(2-ethylhexyl)phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
di-n-octyl phthalate	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
dibenz(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
Surrogate Recovery		Limits									
2-fluorophenol SUR	15 *	21-100	%	1	AJD	10/12/12	5563	10/16/12	23:58	SW3510C8270D	
phenol-D5 SUR	10	10-102	%	1	AJD	10/12/12	5563	10/16/12	23:58	SW3510C8270D	
2,4,6-tribromophenol SUR	58	10-123	%	1	AJD	10/12/12	5563	10/16/12	23:58	SW3510C8270D	
nitrobenzene-D5 SUR	75	35-114	%	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
2-fluorobiphenyl SUR	95	43-116	%	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	
p-terphenyl-D14 SUR	92	33-141	%	1	AJD	10/12/12	5563	10/17/12	0:35	SW3510C8270D	

* The surrogate showed recovery outside the acceptance limits. Matrix interference suspected. No additional sample remains for re-analysis.

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Batch	Analysis			Reference
								Date	Time		
N-nitrosodimethylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
aniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
phenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2-chlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
bis(2-chloroethyl)ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
1,3-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
1,4-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
1,2-dichlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzyl alcohol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
bis(2-chloroisopropyl) ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
hexachloroethane	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
N-nitroso-di-N-propylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
nitrobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
isophorone	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2-nitrophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2,4-dimethylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
bis(2-chloroethoxy)methane	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2,4-dichlorophenol	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
1,2,4-trichlorobenzene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzoic acid	< 50	50	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
4-chloroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
hexachlorobutadiene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4-chloro-3-methylphenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
hexachlorocyclopentadiene	< 10	10	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2,4,6-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2,4,5-trichlorophenol	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2-chloronaphthalene	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
dimethylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2,6-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2,4-dinitrotoluene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
3-nitroaniline	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2,4-dinitrophenol	< 50	50	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
dibenzofuran	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4-nitrophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
diethyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Result	Reporting Limit	Units	Instr Dil'n	Analyst	Prep Date	Batch	Analysis			Reference
								Date	Time		
4-chlorophenyl phenyl ether	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4-nitroaniline	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4,6-dinitro-2-methylphenol	< 20	20	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
azobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
N-nitrosodiphenylamine	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
4-bromophenyl phenyl ether	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
hexachlorobenzene	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
pentachlorophenol	< 10	10	ug/L	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
carbazole	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
di-n-butylphthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
butyl benzyl phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
3,3'-dichlorobenzidine	< 30	30	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
bis(2-ethylhexyl)phthalate	< 5	5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
di-n-octyl phthalate	< 2	2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
dibenz(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
Surrogate Recovery		Limits									
2-fluorophenol SUR	39	21-100	%	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
phenol-D5 SUR	24	10-102	%	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
2,4,6-tribromophenol SUR	110	10-123	%	1	AJD	10/12/12	5563	10/17/12	1:13		SW3510C8270D
nitrobenzene-D5 SUR	78	35-114	%	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
2-fluorobiphenyl SUR	94	43-116	%	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D
p-terphenyl-D14 SUR	94	33-141	%	1	AJD	10/12/12	5563	10/17/12	1:50		SW3510C8270D

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-001

Sample ID: CA-MW-3

Matrix: Water

Sampled: 10/9/12 15:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
dibenzo(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	12:29	MA EPH
Unadjusted C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:02	MA EPH
C9-C18 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:03	MA EPH
C19-C36 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:03	MA EPH
C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:02	MA EPH
Surrogate Recovery										
1-chloro-octadecane SUR	59	40-140	%	1	JLZ	10/12/12	5562	10/16/12	19:03	MA EPH
o-terphenyl SUR	61	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:02	MA EPH
2-fluorobiphenyl SUR	88	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:02	MA EPH
2-bromonaphthalene SUR	85	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:02	MA EPH
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
dibenzo(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:06	MA EPH
Unadjusted C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:24	MA EPH
C9-C18 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:21	MA EPH
C19-C36 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:21	MA EPH
C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:24	MA EPH
Surrogate Recovery										
1-chloro-octadecane SUR	55	40-140	%	1	JLZ	10/12/12	5562	10/16/12	19:21	MA EPH
o-terphenyl SUR	58	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:24	MA EPH
2-fluorobiphenyl SUR	84	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:24	MA EPH
2-bromonaphthalene SUR	81	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:24	MA EPH
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Result	Reporting Limit	Units	Instr Dil'n Factor	Analyst	Prep Date	Batch	Analysis			Reference
								Date	Time		
naphthalene	1.4	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
2-methylnaphthalene	2.9	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
phenanthrene	0.8	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
acenaphthene	0.8	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
fluorene	0.7	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
dibenzo(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	13:43	MA EPH	
Unadjusted C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:47	MA EPH	
C9-C18 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:40	MA EPH	
C19-C36 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:40	MA EPH	
C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	21:47	MA EPH	
Surrogate Recovery											
1-chloro-octadecane SUR	59	40-140	%	1	JLZ	10/12/12	5562	10/16/12	19:40	MA EPH	
o-terphenyl SUR	62	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:47	MA EPH	
2-fluorobiphenyl SUR	84	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:47	MA EPH	
2-bromonaphthalene SUR	87	40-140	%	1	JLZ	10/12/12	5562	10/17/12	21:47	MA EPH	
Limits											

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
naphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
2-methylnaphthalene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
phenanthrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
acenaphthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
acenaphthylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
fluorene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
benzo(a)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
chrysene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
benzo(b)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
benzo(k)fluoranthene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
benzo(a)pyrene	< 0.2	0.2	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
indeno(1,2,3-cd)pyrene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
dibenzo(a,h)anthracene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
benzo(g,h,i)perylene	< 0.5	0.5	ug/L	1	AJD	10/12/12	5562	10/17/12	17:58	MA EPH
Unadjusted C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	22:10	MA EPH
C9-C18 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:58	MA EPH
C19-C36 Aliphatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/16/12	19:58	MA EPH
C11-C22 Aromatics	< 100	100	ug/L	1	JLZ	10/12/12	5562	10/17/12	22:10	MA EPH
Surrogate Recovery										
1-chloro-octadecane SUR	66	40-140	%	1	JLZ	10/12/12	5562	10/16/12	19:58	MA EPH
o-terphenyl SUR	63	40-140	%	1	JLZ	10/12/12	5562	10/17/12	22:10	MA EPH
2-fluorobiphenyl SUR	89	40-140	%	1	JLZ	10/12/12	5562	10/17/12	22:10	MA EPH
2-bromonaphthalene SUR	84	40-140	%	1	JLZ	10/12/12	5562	10/17/12	22:10	MA EPH
Limits										

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-001

Sample ID: CA-MW-3

Matrix: Water

Sampled: 10/9/12 15:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Barium	0.06	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:23	SW7470A
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	20:01	SW3005A6010C

Sample#: 25149-002

Sample ID: CA-MW-6

Matrix: Water

Sampled: 10/9/12 13:20

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Antimony	< 0.006	0.006	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Beryllium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Copper	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:28	SW7470A
Nickel	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Thallium	< 0.002	0.002	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C
Zinc	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:45	SW3005A6010C

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-003

Sample ID: CA-MW-7

Matrix: Water

Sampled: 10/9/12 14:15

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Antimony	< 0.006	0.006	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Beryllium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Copper	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:30	SW7470A
Nickel	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Thallium	< 0.002	0.002	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C
Zinc	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:52	SW3005A6010C

Sample#: 25149-004

Sample ID: CA-MW-8

Matrix: Water

Sampled: 10/9/12 12:12

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Antimony	< 0.006	0.006	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Beryllium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Copper	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:32	SW7470A
Nickel	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Thallium	< 0.002	0.002	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C
Zinc	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:00	SW3005A6010C

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-005

Sample ID: CA-MW-9

Matrix: Water

Sampled: 10/9/12 10:53

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Antimony	< 0.006	0.006	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Beryllium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Copper	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:34	SW7470A
Nickel	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Thallium	< 0.002	0.002	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C
Zinc	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:07	SW3005A6010C

Sample#: 25149-006

Sample ID: CA-MW-10

Matrix: Water

Sampled: 10/9/12 9:42

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Barium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:36	SW7470A
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	20:09	SW3005A6010C

Project ID: MERC 12001151

Job ID: 25149

Sample#: 25149-007

Sample ID: CA-MW-DUP-1

Matrix: Water

Sampled: 10/9/12 0:00

Parameter	Reporting		Instr	Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit					Batch	Date	Time	
Antimony	< 0.006	0.006	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Arsenic	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Beryllium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Cadmium	< 0.004	0.004	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Chromium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Copper	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Lead	< 0.008	0.008	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Mercury	< 0.0002	0.0002	mg/L	1	SAS	10/12/12	5566	10/12/12	18:37	SW7470A
Nickel	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Selenium	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Silver	< 0.007	0.007	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Thallium	< 0.002	0.002	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C
Zinc	< 0.05	0.05	mg/L	1	SAS	10/12/12	5569	10/15/12	21:15	SW3005A6010C

Quality Control Report



124 Heritage Avenue Unit 16
Portsmouth, NH 03801
www.absoluteresourceassociates.com



Case Narrative
Lab # 25149

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 4 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

ABN: Sample CA-MW-10 did not meet acceptance criteria for the surrogate 2-fluorophenol due suspected matrix interference No additional sample remains for re-analysis.

Laboratory Control Sample Results

ABN: The LCS/D5530 did not meet the acceptance criteria for N-nitrosodimethylamine, phenol, hexachlorocyclopentadiene, dimethylphthalate, 4-nitrophenol. Since <10% of the compounds were outside of the acceptance criteria, reanalysis is not required.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

PAH: Dilution was required due to hydrocarbon interferences for the following samples: CA-MW-8.

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Association Table -

Analysis		QC Number	Field ID	Lab ID
EPH in water by	MA EPH			
		5562	CA-MW-3 CA-MW-6 CA-MW-7 CA-MW-DUP-1	25149-001 25149-002 25149-003 25149-007
VPH in water by MA	MA VPH			
		1202475	CA-MW-6 CA-MW-7 CA-MW-DUP-1 Trip Blank	25149-002 25149-003 25149-007 25149-009
Antimony in water by	SW3005A6010C			
		5569	CA-MW-3 CA-MW-6 CA-MW-7 CA-MW-8 CA-MW-9 CA-MW-10 CA-MW-DUP-1	25149-001 25149-002 25149-003 25149-004 25149-005 25149-006 25149-007
Acid & Base/Neutral	SW3510C8270D			
		5563	CA-MW-6 CA-MW-7 CA-MW-8 CA-MW-9 CA-MW-10 CA-MW-DUP-1	25149-002 25149-003 25149-004 25149-005 25149-006 25149-007
VOCs in water by 8260	SW5030B8260B			
		1202467	CA-MW-6 CA-MW-7 CA-MW-9 CA-MW-10 CA-MW-DUP-1 Trip Blank	25149-002 25149-003 25149-005 25149-006 25149-007 25149-008
Mercury in water by 7470	SW7470A			
		5566	CA-MW-3 CA-MW-6 CA-MW-7 CA-MW-8 CA-MW-9 CA-MW-10 CA-MW-DUP-1	25149-001 25149-002 25149-003 25149-004 25149-005 25149-006 25149-007

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA VPH	BLK1202475	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		<	100	ug/L				
		Unadjusted C9-C12 Aliphatics		<	100	ug/L				
		methyl t-butyl ether (MTBE)		<	2	ug/L				
		benzene		<	1	ug/L				
		toluene		<	2	ug/L				
		ethylbenzene		<	2	ug/L				
		m&p-xlenes		<	2	ug/L				
		o-xylene		<	2	ug/L				
		naphthalene		<	5	ug/L				
		C9-C10 Aromatics		<	100	ug/L				
		2,5-dibromotoluene as		113	%			70 130		
		2,5-dibromotoluene as		114	%			70 130		
MA VPH	LCS1202475	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		820	ug/L	750	109	70 130		
		Unadjusted C9-C12 Aliphatics		520	ug/L	550	94	70 130		
		methyl t-butyl ether (MTBE)		170	ug/L	150	111	70 130		
		benzene		50	ug/L	50	101	70 130		
		toluene		160	ug/L	150	103	70 130		
		ethylbenzene		49	ug/L	50	99	70 130		
		m&p-xlenes		200	ug/L	200	102	70 130		
		o-xylene		100	ug/L	100	103	70 130		
		naphthalene		100	ug/L	100	101	70 130		
		C9-C10 Aromatics		100	ug/L	100	100	70 130		
		2,5-dibromotoluene as		370	%			70 130		
		2,5-dibromotoluene as		361	%			70 130		
MA VPH	LCSD1202475	C5-C8 Aliphatics								
		Unadjusted C5-C8 Aliphatics		800	ug/L	750	107	70 130	2	25
		Unadjusted C9-C12 Aliphatics		520	ug/L	550	94	70 130	0	25
		methyl t-butyl ether (MTBE)		170	ug/L	150	112	70 130	2	25
		benzene		51	ug/L	50	102	70 130	1	25
		toluene		160	ug/L	150	105	70 130	1	25
		ethylbenzene		48	ug/L	50	96	70 130	3	25
		m&p-xlenes		210	ug/L	200	103	70 130	1	25
		o-xylene		100	ug/L	100	104	70 130	1	25
		naphthalene		100	ug/L	100	101	70 130	1	25
		C9-C10 Aromatics		100	ug/L	100	101	70 130	1	25
		2,5-dibromotoluene as		375	%			70 130		
		2,5-dibromotoluene as		364	%			70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	BLK1202467	dichlorodifluoromethane		<	2	ug/L				
		chloromethane		<	2	ug/L				
		vinyl chloride		<	2	ug/L				
		bromomethane		<	2	ug/L				
		chloroethane		<	2	ug/L				
		trichlorofluoromethane		<	2	ug/L				
		diethyl ether		<	10	ug/L				
		acetone		<	50	ug/L				
		1,1-dichloroethene		<	1	ug/L				
		methylene chloride		<	5	ug/L				
		carbon disulfide		<	2	ug/L				
		methyl t-butyl ether (MTBE)		<	2	ug/L				
		trans-1,2-dichloroethene		<	2	ug/L				
		1,1-dichloroethane		<	2	ug/L				
		2-butanone (MEK)		<	10	ug/L				
		2,2-dichloropropane		<	2	ug/L				
		cis-1,2-dichloroethene		<	2	ug/L				
		chloroform		<	2	ug/L				
		bromochloromethane		<	2	ug/L				
		tetrahydrofuran (THF)		<	10	ug/L				
		1,1,1-trichloroethane		<	2	ug/L				
		1,1-dichloropropene		<	2	ug/L				
		carbon tetrachloride		<	2	ug/L				
		1,2-dichloroethane		<	2	ug/L				
		benzene		<	2	ug/L				
		trichloroethene		<	2	ug/L				
		1,2-dichloropropane		<	2	ug/L				
		bromodichloromethane		<	0.6	ug/L				
		dibromomethane		<	2	ug/L				
		4-methyl-2-pentanone (MIBK)		<	10	ug/L				
		cis-1,3-dichloropropene		<	2	ug/L				
		toluene		<	2	ug/L				
		trans-1,3-dichloropropene		<	2	ug/L				
		2-hexanone		<	10	ug/L				
		1,1,2-trichloroethane		<	2	ug/L				
		1,3-dichloropropane		<	2	ug/L				
		tetrachloroethene		<	2	ug/L				
		dibromochloromethane		<	2	ug/L				
		1,2-dibromoethane (EDB)		<	2	ug/L				
		chlorobenzene		<	2	ug/L				
		1,1,1,2-tetrachloroethane		<	2	ug/L				
		ethylbenzene		<	2	ug/L				
		m&p-xlenes		<	2	ug/L				
		o-xylene		<	2	ug/L				
		styrene		<	2	ug/L				
		bromoform		<	2	ug/L				
		isopropylbenzene		<	2	ug/L				
		1,1,2,2-tetrachloroethane		<	2	ug/L				
		1,2,3-trichloropropane		<	2	ug/L				
		n-propylbenzene		<	2	ug/L				
		bromobenzene		<	2	ug/L				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	BLK1202467	1,3,5-trimethylbenzene		<	2	ug/L				
		2-chlorotoluene		<	2	ug/L				
		4-chlorotoluene		<	2	ug/L				
		tert-butylbenzene		<	2	ug/L				
		1,2,4-trimethylbenzene		<	2	ug/L				
		sec-butylbenzene		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		4-isopropyltoluene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		n-butylbenzene		<	2	ug/L				
		1,2-dibromo-3-chloropropane		<	2	ug/L				
		1,2,4-trichlorobenzene		<	2	ug/L				
		hexachlorobutadiene		<	0.5	ug/L				
		naphthalene		<	5	ug/L				
		1,2,3-trichlorobenzene		<	2	ug/L				
		dibromofluoromethane SUR		102	%		78	114		
		toluene-D8 SUR		101	%		88	110		
		4-bromofluorobenzene SUR		107	%		86	115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1202467	dichlorodifluoromethane		18	ug/L	20	89	70 130		
		chloromethane		19	ug/L	20	97	70 130		
		vinyl chloride		19	ug/L	20	97	70 130		
		bromomethane		15	ug/L	20	73	70 130		
		chloroethane		21	ug/L	20	106	70 130		
		trichlorofluoromethane		19	ug/L	20	96	70 130		
		diethyl ether		20	ug/L	20	102	70 130		
		acetone	<	50	ug/L	20	109			
		1,1-dichloroethene		19	ug/L	20	93	70 130		
		methylene chloride		22	ug/L	20	108	70 130		
		carbon disulfide		24	ug/L	20	122	70 130		
		methyl t-butyl ether (MTBE)		20	ug/L	20	100	70 130		
		trans-1,2-dichloroethene		20	ug/L	20	101	70 130		
		1,1-dichloroethane		20	ug/L	20	99	70 130		
		2-butanone (MEK)		23	ug/L	20	114	70 130		
		2,2-dichloropropane		19	ug/L	20	95	70 130		
		cis-1,2-dichloroethene		21	ug/L	20	103	70 130		
		chloroform		20	ug/L	20	101	70 130		
		bromochloromethane		20	ug/L	20	102	70 130		
		tetrahydrofuran (THF)		20	ug/L	20	101	70 130		
		1,1,1-trichloroethane		20	ug/L	20	101	70 130		
		1,1-dichloropropene		21	ug/L	20	105	70 130		
		carbon tetrachloride		19	ug/L	20	96	70 130		
		1,2-dichloroethane		19	ug/L	20	97	70 130		
		benzene		20	ug/L	20	99	70 130		
		trichloroethene		18	ug/L	20	92	70 130		
		1,2-dichloropropane		19	ug/L	20	97	70 130		
		bromodichloromethane		20	ug/L	20	100	70 130		
		dibromomethane		20	ug/L	20	101	70 130		
		4-methyl-2-pentanone (MIBK)		22	ug/L	20	109	70 130		
		cis-1,3-dichloropropene		20	ug/L	20	101	70 130		
		toluene		22	ug/L	20	108	70 130		
		trans-1,3-dichloropropene		19	ug/L	20	97	70 130		
		2-hexanone		24	ug/L	20	118	70 130		
		1,1,2-trichloroethane		20	ug/L	20	98	70 130		
		1,3-dichloropropane		19	ug/L	20	96	70 130		
		tetrachloroethene		19	ug/L	20	94	70 130		
		dibromochloromethane		19	ug/L	20	95	70 130		
		1,2-dibromoethane (EDB)		20	ug/L	20	98	70 130		
		chlorobenzene		21	ug/L	20	103	70 130		
		1,1,1,2-tetrachloroethane		19	ug/L	20	96	70 130		
		ethylbenzene		19	ug/L	20	97	70 130		
		m&p-xlenes		41	ug/L	40	102	70 130		
		o-xylene		20	ug/L	20	102	70 130		
		styrene		20	ug/L	20	102	70 130		
		bromoform		20	ug/L	20	98	70 130		
		isopropylbenzene		19	ug/L	20	96	70 130		
		1,1,2,2-tetrachloroethane		20	ug/L	20	101	70 130		
		1,2,3-trichloropropane		19	ug/L	20	93	70 130		
		n-propylbenzene		19	ug/L	20	94	70 130		
		bromobenzene		19	ug/L	20	95	70 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCS1202467	1,3,5-trimethylbenzene		19	ug/L	20	97	70 130		
		2-chlorotoluene		19	ug/L	20	97	70 130		
		4-chlorotoluene		20	ug/L	20	98	70 130		
		tert-butylbenzene		19	ug/L	20	95	70 130		
		1,2,4-trimethylbenzene		19	ug/L	20	95	70 130		
		sec-butylbenzene		20	ug/L	20	98	70 130		
		1,3-dichlorobenzene		19	ug/L	20	96	70 130		
		4-isopropyltoluene		19	ug/L	20	97	70 130		
		1,4-dichlorobenzene		19	ug/L	20	95	70 130		
		1,2-dichlorobenzene		20	ug/L	20	98	70 130		
		n-butylbenzene		19	ug/L	20	96	70 130		
		1,2-dibromo-3-chloropropane		21	ug/L	20	103	70 130		
		1,2,4-trichlorobenzene		18	ug/L	20	89	70 130		
		hexachlorobutadiene		18	ug/L	20	88	70 130		
		naphthalene		20	ug/L	20	102	70 130		
		1,2,3-trichlorobenzene		19	ug/L	20	96	70 130		
		dibromofluoromethane SUR		102	%			78 114		
		toluene-D8 SUR		103	%			88 110		
		4-bromofluorobenzene SUR		105	%			86 115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCSD1202467	dichlorodifluoromethane		17	ug/L	20	87	70 130	2	20
		chloromethane		17	ug/L	20	83	70 130	16	20
		vinyl chloride		18	ug/L	20	91	70 130	6	20
		bromomethane		14	ug/L	20	72	70 130	2	20
		chloroethane		20	ug/L	20	99	70 130	7	20
		trichlorofluoromethane		19	ug/L	20	96	70 130	0	20
		diethyl ether		19	ug/L	20	97	70 130	6	20
		acetone	<	50	ug/L	20	105		4	20
		1,1-dichloroethene		17	ug/L	20	87	70 130	7	20
		methylene chloride		20	ug/L	20	101	70 130	7	20
		carbon disulfide		23	ug/L	20	114	70 130	6	20
		methyl t-butyl ether (MTBE)		19	ug/L	20	95	70 130	5	20
		trans-1,2-dichloroethene		19	ug/L	20	94	70 130	7	20
		1,1-dichloroethane		18	ug/L	20	91	70 130	9	20
		2-butanone (MEK)		23	ug/L	20	114	70 130	0	20
		2,2-dichloropropane		17	ug/L	20	87	70 130	9	20
		cis-1,2-dichloroethene		19	ug/L	20	95	70 130	9	20
		chloroform		19	ug/L	20	94	70 130	7	20
		bromochloromethane		19	ug/L	20	96	70 130	6	20
		tetrahydrofuran (THF)		19	ug/L	20	96	70 130	5	20
		1,1,1-trichloroethane		19	ug/L	20	93	70 130	8	20
		1,1-dichloropropene		20	ug/L	20	99	70 130	7	20
		carbon tetrachloride		18	ug/L	20	92	70 130	4	20
		1,2-dichloroethane		18	ug/L	20	90	70 130	8	20
		benzene		18	ug/L	20	91	70 130	8	20
		trichloroethene		17	ug/L	20	85	70 130	8	20
		1,2-dichloropropane		18	ug/L	20	88	70 130	9	20
		bromodichloromethane		19	ug/L	20	93	70 130	8	20
		dibromomethane		18	ug/L	20	92	70 130	9	20
		4-methyl-2-pentanone (MIBK)		22	ug/L	20	108	70 130	1	20
		cis-1,3-dichloropropene		19	ug/L	20	93	70 130	7	20
		toluene		19	ug/L	20	97	70 130	11	20
		trans-1,3-dichloropropene		18	ug/L	20	91	70 130	6	20
		2-hexanone		23	ug/L	20	116	70 130	2	20
		1,1,2-trichloroethane		19	ug/L	20	93	70 130	5	20
		1,3-dichloropropane		19	ug/L	20	94	70 130	2	20
		tetrachloroethene		19	ug/L	20	93	70 130	1	20
		dibromochloromethane		19	ug/L	20	93	70 130	2	20
		1,2-dibromoethane (EDB)		19	ug/L	20	95	70 130	3	20
		chlorobenzene		19	ug/L	20	96	70 130	7	20
		1,1,1,2-tetrachloroethane		19	ug/L	20	95	70 130	2	20
		ethylbenzene		19	ug/L	20	94	70 130	4	20
		m&p-xlenes		38	ug/L	40	96	70 130	6	20
		o-xylene		19	ug/L	20	96	70 130	6	20
		styrene		20	ug/L	20	98	70 130	4	20
		bromoform		20	ug/L	20	98	70 130	0	20
		isopropylbenzene		18	ug/L	20	92	70 130	4	20
		1,1,2,2-tetrachloroethane		20	ug/L	20	101	70 130	0	20
		1,2,3-trichloropropane		19	ug/L	20	95	70 130	2	20
		n-propylbenzene		18	ug/L	20	91	70 130	3	20
		bromobenzene		18	ug/L	20	92	70 130	3	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW5030B8260B	LCSD1202467	1,3,5-trimethylbenzene		18	ug/L	20	92	70 130	5	20
		2-chlorotoluene		18	ug/L	20	89	70 130	8	20
		4-chlorotoluene		19	ug/L	20	95	70 130	3	20
		tert-butylbenzene		18	ug/L	20	90	70 130	6	20
		1,2,4-trimethylbenzene		18	ug/L	20	92	70 130	4	20
		sec-butylbenzene		19	ug/L	20	96	70 130	2	20
		1,3-dichlorobenzene		18	ug/L	20	91	70 130	6	20
		4-isopropyltoluene		18	ug/L	20	92	70 130	6	20
		1,4-dichlorobenzene		18	ug/L	20	91	70 130	4	20
		1,2-dichlorobenzene		18	ug/L	20	92	70 130	6	20
		n-butylbenzene		19	ug/L	20	93	70 130	3	20
		1,2-dibromo-3-chloropropane		21	ug/L	20	104	70 130	1	20
		1,2,4-trichlorobenzene		17	ug/L	20	87	70 130	2	20
		hexachlorobutadiene		18	ug/L	20	88	70 130	0	20
		naphthalene		20	ug/L	20	100	70 130	1	20
		1,2,3-trichlorobenzene		19	ug/L	20	94	70 130	2	20
		dibromofluoromethane SUR		102	%			78 114		
		toluene-D8 SUR		101	%			88 110		
		4-bromofluorobenzene SUR		106	%			86 115		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	BLK5562	naphthalene		<	ug/L	0.5				
		2-methylnaphthalene		<	ug/L	0.5				
		phenanthrene		<	ug/L	0.5				
		acenaphthene		<	ug/L	0.5				
		acenaphthylene		<	ug/L	0.5				
		fluorene		<	ug/L	0.5				
		anthracene		<	ug/L	0.5				
		fluoranthene		<	ug/L	0.5				
		pyrene		<	ug/L	0.5				
		benzo(a)anthracene		<	ug/L	0.5				
		chrysene		<	ug/L	0.5				
		benzo(b)fluoranthene		<	ug/L	0.5				
		benzo(k)fluoranthene		<	ug/L	0.5				
		benzo(a)pyrene		<	ug/L	0.2				
		indeno(1,2,3-cd)pyrene		<	ug/L	0.5				
		dibenzo(a,h)anthracene		<	ug/L	0.5				
		benzo(g,h,i)perylene		<	ug/L	0.5				
		Unadjusted C11-C22 Aromatics		<	ug/L	100				
		C9-C18 Aliphatics		<	ug/L	100				
		C19-C36 Aliphatics		<	ug/L	100				
		C11-C22 Aromatics		<	ug/L	100				
		1-chloro-octadecane SUR		65	%			40	140	
		o-terphenyl SUR		85	%			40	140	
		2-fluorobiphenyl SUR		126	%			40	140	
		2-bromonaphthalene SUR		135	%			40	140	
MA EPH	LCS5562	naphthalene		40	ug/L	60	67	40	140	
		2-methylnaphthalene		39	ug/L	60	64	40	140	
		phenanthrene		47	ug/L	60	78	40	140	
		acenaphthene		43	ug/L	60	72	40	140	
		acenaphthylene		45	ug/L	60	75	40	140	
		fluorene		38	ug/L	60	63	40	140	
		anthracene		43	ug/L	60	71	40	140	
		fluoranthene		46	ug/L	60	76	40	140	
		pyrene		54	ug/L	60	91	40	140	
		benzo(a)anthracene		49	ug/L	60	82	40	140	
		chrysene		51	ug/L	60	86	40	140	
		benzo(b)fluoranthene		51	ug/L	60	84	40	140	
		benzo(k)fluoranthene		55	ug/L	60	92	40	140	
		benzo(a)pyrene		51	ug/L	60	84	40	140	
		indeno(1,2,3-cd)pyrene		51	ug/L	60	85	40	140	
		dibenzo(a,h)anthracene		52	ug/L	60	86	40	140	
		benzo(g,h,i)perylene		51	ug/L	60	85	40	140	
		Unadjusted C11-C22 Aromatics		760	ug/L	1020	75	40	140	
		C9-C18 Aliphatics		280	ug/L	360	78	40	140	
		C19-C36 Aliphatics		430	ug/L	480	90	40	140	
		C11-C22 Aromatics		<	ug/L	100				
		1-chloro-octadecane SUR		67	%			40	140	
		o-terphenyl SUR		85	%			40	140	
		2-fluorobiphenyl SUR		121	%			40	140	
		2-bromonaphthalene SUR		113	%			40	140	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
MA EPH	LCSD5562	naphthalene		43	ug/L	60	71	40 140	7	25
		2-methylnaphthalene		41	ug/L	60	68	40 140	5	25
		phenanthrene		46	ug/L	60	77	40 140	1	25
		acenaphthene		44	ug/L	60	74	40 140	2	25
		acenaphthylene		46	ug/L	60	76	40 140	1	25
		fluorene		41	ug/L	60	69	40 140	9	25
		anthracene		42	ug/L	60	71	40 140	1	25
		fluoranthene		38	ug/L	60	64	40 140	18	25
		pyrene		53	ug/L	60	88	40 140	3	25
		benzo(a)anthracene		51	ug/L	60	84	40 140	3	25
		chrysene		52	ug/L	60	87	40 140	1	25
		benzo(b)fluoranthene		54	ug/L	60	89	40 140	6	25
		benzo(k)fluoranthene		52	ug/L	60	87	40 140	5	25
		benzo(a)pyrene		51	ug/L	60	85	40 140	0	25
		indeno(1,2,3-cd)pyrene		51	ug/L	60	85	40 140	0	25
		dibenzo(a,h)anthracene		52	ug/L	60	87	40 140	0	25
		benzo(g,h,i)perylene		51	ug/L	60	85	40 140	0	25
		Unadjusted C11-C22 Aromatics		680	ug/L	1020	67	40 140	11	25
		C9-C18 Aliphatics		290	ug/L	360	81	40 140	3	25
		C19-C36 Aliphatics		490	ug/L	480	102	40 140	12	25
		C11-C22 Aromatics	<	100	ug/L					
		1-chloro-octadecane SUR		71	%			40 140		
		o-terphenyl SUR		79	%			40 140		
		2-fluorobiphenyl SUR		113	%			40 140		
		2-bromonaphthalene SUR		103	%			40 140		

Lab Number: NBT LCS5562
Sample Designation: Laboratory Control Sample 5562W
Date Sampled: N/A
Date Received: N/A
Date Extracted: 10/11/12
Matrix: Water
Containers: N/A
Sample Preservation: N/A
Temperature: N/A
Dilution Factor: 1
Analyst: JLZ

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

	LCS Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.3%	<5.0%	10/17/12
2-methylnaphthalene	0.3%	<5.0%	10/17/12

Lab Number: NBT LCSD5562
Sample Designation: Laboratory Control Sample Duplicate 5562W
Date Sampled: N/A
Date Received: N/A
Date Extracted: 10/11/12
Matrix: Water
Containers: N/A
Sample Preservation: N/A
Temperature: N/A
Dilution Factor: 1
Analyst: JLZ

NAPHTHALENE BREAKTHROUGH CALCULATION

Method for Ranges: MADEP EPH 2004-1.1

Method for Target Analytes: EPA 8270C

	LCSD Aliphatic Breakthrough (%)	Acceptance Criteria	Date of Analysis
naphthalene	0.2%	<5.0%	10/17/12
2-methylnaphthalene	0.3%	<5.0%	10/17/12

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	BLK5563	N-nitrosodimethylamine		<	2	ug/L				
		aniline		<	2	ug/L				
		phenol		<	2	ug/L				
		2-chlorophenol		<	5	ug/L				
		bis(2-chloroethyl)ether		<	2	ug/L				
		1,3-dichlorobenzene		<	2	ug/L				
		1,4-dichlorobenzene		<	2	ug/L				
		1,2-dichlorobenzene		<	2	ug/L				
		benzyl alcohol		<	2	ug/L				
		2-methylphenol		<	2	ug/L				
		bis(2-chloroisopropyl) ether		<	2	ug/L				
		hexachloroethane		<	2	ug/L				
		N-nitroso-di-N-propylamine		<	2	ug/L				
		4-methylphenol		<	2	ug/L				
		nitrobenzene		<	2	ug/L				
		isophorone		<	5	ug/L				
		2-nitrophenol		<	2	ug/L				
		2,4-dimethylphenol		<	2	ug/L				
		bis(2-chloroethoxy)methane		<	5	ug/L				
		2,4-dichlorophenol		<	5	ug/L				
		1,2,4-trichlorobenzene		<	5	ug/L				
		naphthalene		<	0.5	ug/L				
		benzoic acid		<	50	ug/L				
		4-chloroaniline		<	2	ug/L				
		hexachlorobutadiene		<	2	ug/L				
		4-chloro-3-methylphenol		<	2	ug/L				
		2-methylnaphthalene		<	0.5	ug/L				
		hexachlorocyclopentadiene		<	10	ug/L				
		2,4,6-trichlorophenol		<	2	ug/L				
		2,4,5-trichlorophenol		<	2	ug/L				
		2-chloronaphthalene		<	5	ug/L				
		2-nitroaniline		<	2	ug/L				
		acenaphthylene		<	0.5	ug/L				
		dimethylphthalate		<	5	ug/L				
		2,6-dinitrotoluene		<	2	ug/L				
		2,4-dinitrotoluene		<	2	ug/L				
		acenaphthene		<	0.5	ug/L				
		3-nitroaniline		<	2	ug/L				
		2,4-dinitrophenol		<	50	ug/L				
		dibenzofuran		<	0.5	ug/L				
		4-nitrophenol		<	10	ug/L				
		fluorene		<	0.5	ug/L				
		diethyl phthalate		<	5	ug/L				
		4-chlorophenyl phenyl ether		<	5	ug/L				
		4-nitroaniline		<	5	ug/L				
		4,6-dinitro-2-methylphenol		<	20	ug/L				
		azobenzene		<	2	ug/L				
		N-nitrosodiphenylamine		<	2	ug/L				
		4-bromophenyl phenyl ether		<	2	ug/L				
		hexachlorobenzene		<	2	ug/L				
		pentachlorophenol		<	10	ug/L				

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	BLK5563	phenanthrene		<	0.5	ug/L				
		anthracene		<	0.5	ug/L				
		carbazole		<	2	ug/L				
		di-n-butylphthalate		<	5	ug/L				
		fluoranthene		<	0.5	ug/L				
		benzidine		<	30	ug/L				
		pyrene		<	0.5	ug/L				
		butyl benzyl phthalate		<	5	ug/L				
		benzo(a)anthracene		<	0.5	ug/L				
		chrysene		<	0.5	ug/L				
		3,3'-dichlorobenzidine		<	30	ug/L				
		bis(2-ethylhexyl)phthalate		<	5	ug/L				
		di-n-octyl phthalate		<	2	ug/L				
		benzo(b)fluoranthene		<	0.5	ug/L				
		benzo(k)fluoranthene		<	0.5	ug/L				
		benzo(a)pyrene		<	0.2	ug/L				
		indeno(1,2,3-cd)pyrene		<	0.5	ug/L				
		dibenzo(a,h)anthracene		<	0.5	ug/L				
		benzo(g,h,i)perylene		<	0.5	ug/L				
		2-fluorophenol SUR		28	%			21	100	
		phenol-D5 SUR		17	%			10	102	
		2,4,6-tribromophenol SUR		68	%			10	123	
		nitrobenzene-D5 SUR		57	%			35	114	
		2-fluorobiphenyl SUR		67	%			43	116	
		p-terphenyl-D14 SUR		87	%			33	141	

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCS5563	N-nitrosodimethylamine		9	ug/L	40	22	* 40 140		
		aniline		17	ug/L	40	42	40 140		
		phenol		6	ug/L	40	16	* 30 130		
		2-chlorophenol		22	ug/L	40	54	30 130		
		bis(2-chloroethyl)ether		26	ug/L	40	66	40 140		
		1,3-dichlorobenzene		17	ug/L	40	43	40 140		
		1,4-dichlorobenzene		18	ug/L	40	46	40 140		
		1,2-dichlorobenzene		20	ug/L	40	49	40 140		
		benzyl alcohol		16	ug/L	40	41	30 130		
		2-methylphenol		17	ug/L	40	41	30 130		
		bis(2-chloroisopropyl) ether		23	ug/L	40	57	40 140		
		hexachloroethane		18	ug/L	40	44	40 140		
		N-nitroso-di-N-propylamine		22	ug/L	40	56	40 140		
		4-methylphenol		16	ug/L	40	39	30 130		
		nitrobenzene		22	ug/L	40	55	40 140		
		isophorone		24	ug/L	40	59	40 140		
		2-nitrophenol		25	ug/L	40	62	30 130		
		2,4-dimethylphenol		12	ug/L	40	30	30 130		
		bis(2-chloroethoxy)methane		25	ug/L	40	62	40 140		
		2,4-dichlorophenol		25	ug/L	40	64	30 130		
		1,2,4-trichlorobenzene		21	ug/L	40	51	40 140		
		naphthalene		22	ug/L	40	55	40 140		
		benzoic acid	<	50	ug/L					
		4-chloroaniline		29	ug/L	40	71	40 140		
		hexachlorobutadiene		20	ug/L	40	49	40 140		
		4-chloro-3-methylphenol		23	ug/L	40	57	30 130		
		2-methylnaphthalene		22	ug/L	40	56	40 140		
		hexachlorocyclopentadiene	<	10	ug/L	40	8	* 40 140		
		2,4,6-trichlorophenol		29	ug/L	40	72	30 130		
		2,4,5-trichlorophenol		26	ug/L	40	64	30 130		
		2-chloronaphthalene		24	ug/L	40	60	40 140		
		2-nitroaniline		27	ug/L	40	68	40 140		
		acenaphthylene		24	ug/L	40	59	40 140		
		dimethylphthalate		12	ug/L	40	30	* 40 140		
		2,6-dinitrotoluene		28	ug/L	40	71	40 140		
		2,4-dinitrotoluene		27	ug/L	40	68	40 140		
		acenaphthene		25	ug/L	40	61	40 140		
		3-nitroaniline		31	ug/L	40	78	40 140		
		2,4-dinitrophenol	<	50	ug/L					
		dibenzofuran		24	ug/L	40	60	40 140		
		4-nitrophenol	<	10	ug/L	40	19	* 30 130		
		fluorene		26	ug/L	40	65	40 140		
		diethyl phthalate		21	ug/L	40	52	40 140		
		4-chlorophenyl phenyl ether		26	ug/L	40	65	40 140		
		4-nitroaniline		24	ug/L	40	61	40 140		
		4,6-dinitro-2-methylphenol	<	20	ug/L					
		azobenzene		24	ug/L	40	60	40 140		
		N-nitrosodiphenylamine		27	ug/L	40	68	40 140		
		4-bromophenyl phenyl ether		25	ug/L	40	63	40 140		
		hexachlorobenzene		25	ug/L	40	61	40 140		
		pentachlorophenol		25	ug/L	40	62	30 130		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCS5563	phenanthrene		25	ug/L	40	63	40 140		
		anthracene		24	ug/L	40	60	40 140		
		carbazole		27	ug/L	40	66	40 140		
		di-n-butylphthalate		26	ug/L	40	64	40 140		
		fluoranthene		26	ug/L	40	64	40 140		
		benzidine		<	30	ug/L				
		pyrene		32	ug/L	40	80	40 140		
		butyl benzyl phthalate		25	ug/L	40	62	40 140		
		benzo(a)anthracene		27	ug/L	40	67	40 140		
		chrysene		26	ug/L	40	64	40 140		
		3,3'-dichlorobenzidine		<	30	ug/L				
		bis(2-ethylhexyl)phthalate		28	ug/L	40	69	40 140		
		di-n-octyl phthalate		35	ug/L	40	86	40 140		
		benzo(b)fluoranthene		30	ug/L	40	74	40 140		
		benzo(k)fluoranthene		34	ug/L	40	85	40 140		
		benzo(a)pyrene		30	ug/L	40	75	40 140		
		indeno(1,2,3-cd)pyrene		31	ug/L	40	77	40 140		
		dibenzo(a,h)anthracene		31	ug/L	40	78	40 140		
		benzo(g,h,i)perylene		31	ug/L	40	78	40 140		
		2-fluorophenol SUR		28	%			21 100		
		phenol-D5 SUR		16	%			10 102		
		2,4,6-tribromophenol SUR		75	%			10 123		
		nitrobenzene-D5 SUR		57	%			35 114		
		2-fluorobiphenyl SUR		67	%			43 116		
		p-terphenyl-D14 SUR		82	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCSD5563	N-nitrosodimethylamine		8	ug/L	40	20	* 40 140	6	20
		aniline		17	ug/L	40	42	40 140	1	20
		phenol		7	ug/L	40	17	* 30 130	6	20
		2-chlorophenol		23	ug/L	40	57	30 130	6	20
		bis(2-chloroethyl)ether		28	ug/L	40	69	40 140	5	20
		1,3-dichlorobenzene		17	ug/L	40	43	40 140	1	20
		1,4-dichlorobenzene		18	ug/L	40	46	40 140	1	20
		1,2-dichlorobenzene		20	ug/L	40	49	40 140	0	20
		benzyl alcohol		16	ug/L	40	40	30 130	2	20
		2-methylphenol		17	ug/L	40	43	30 130	4	20
		bis(2-chloroisopropyl) ether		23	ug/L	40	57	40 140	0	20
		hexachloroethane		16	ug/L	40	40	40 140	9	20
		N-nitroso-di-N-propylamine		23	ug/L	40	57	40 140	1	20
		4-methylphenol		17	ug/L	40	42	30 130	7	20
		nitrobenzene		23	ug/L	40	57	40 140	4	20
		isophorone		24	ug/L	40	59	40 140	0	20
		2-nitrophenol		26	ug/L	40	64	30 130	3	20
		2,4-dimethylphenol		14	ug/L	40	34	30 130	14	20
		bis(2-chloroethoxy)methane		25	ug/L	40	63	40 140	1	20
		2,4-dichlorophenol		27	ug/L	40	68	30 130	7	20
		1,2,4-trichlorobenzene		21	ug/L	40	53	40 140	3	20
		naphthalene		22	ug/L	40	56	40 140	1	20
		benzoic acid	<	50	ug/L					
		4-chloroaniline		28	ug/L	40	70	40 140	1	20
		hexachlorobutadiene		20	ug/L	40	49	40 140	1	20
		4-chloro-3-methylphenol		25	ug/L	40	63	30 130	10	20
		2-methylnaphthalene		22	ug/L	40	56	40 140	1	20
		hexachlorocyclopentadiene	<	10	ug/L	40	8	* 40 140	1	20
		2,4,6-trichlorophenol		30	ug/L	40	76	30 130	6	20
		2,4,5-trichlorophenol		29	ug/L	40	72	30 130	11	20
		2-chloronaphthalene		24	ug/L	40	59	40 140	1	20
		2-nitroaniline		27	ug/L	40	68	40 140	1	20
		acenaphthylene		23	ug/L	40	58	40 140	3	20
		dimethylphthalate		11	ug/L	40	28	* 40 140	7	20
		2,6-dinitrotoluene		28	ug/L	40	70	40 140	1	20
		2,4-dinitrotoluene		27	ug/L	40	67	40 140	1	20
		acenaphthene		23	ug/L	40	59	40 140	5	20
		3-nitroaniline		31	ug/L	40	78	40 140	0	20
		2,4-dinitrophenol	<	50	ug/L					
		dibenzofuran		24	ug/L	40	59	40 140	2	20
		4-nitrophenol	<	10	ug/L	40	21	* 30 130	13	20
		fluorene		26	ug/L	40	64	40 140	1	20
		diethyl phthalate		21	ug/L	40	54	40 140	3	20
		4-chlorophenyl phenyl ether		26	ug/L	40	65	40 140	1	20
		4-nitroaniline		25	ug/L	40	62	40 140	1	20
		4,6-dinitro-2-methylphenol		21	ug/L					
		azobenzene		24	ug/L	40	60	40 140	0	20
		N-nitrosodiphenylamine		25	ug/L	40	62	40 140	9	20
		4-bromophenyl phenyl ether		24	ug/L	40	61	40 140	3	20
		hexachlorobenzene		24	ug/L	40	60	40 140	2	20
		pentachlorophenol		28	ug/L	40	71	30 130	13	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3510C8270D	LCSD5563	phenanthrene		25	ug/L	40	62	40 140	3	20
		anthracene		23	ug/L	40	58	40 140	4	20
		carbazole		26	ug/L	40	64	40 140	3	20
		di-n-butylphthalate		25	ug/L	40	62	40 140	4	20
		fluoranthene		24	ug/L	40	61	40 140	5	20
		benzidine	<	30	ug/L					
		pyrene		29	ug/L	40	74	40 140	9	20
		butyl benzyl phthalate		24	ug/L	40	59	40 140	5	20
		benzo(a)anthracene		25	ug/L	40	63	40 140	6	20
		chrysene		24	ug/L	40	61	40 140	6	20
		3,3'-dichlorobenzidine	<	30	ug/L					
		bis(2-ethylhexyl)phthalate		26	ug/L	40	64	40 140	7	20
		di-n-octyl phthalate		32	ug/L	40	81	40 140	6	20
		benzo(b)fluoranthene		31	ug/L	40	77	40 140	4	20
		benzo(k)fluoranthene		28	ug/L	40	71	40 140	18	20
		benzo(a)pyrene		28	ug/L	40	71	40 140	6	20
		indeno(1,2,3-cd)pyrene		29	ug/L	40	72	40 140	7	20
		dibenzo(a,h)anthracene		29	ug/L	40	72	40 140	7	20
		benzo(g,h,i)perylene		29	ug/L	40	72	40 140	8	20
		2-fluorophenol SUR		29	%			21 100		
		phenol-D5 SUR		17	%			10 102		
		2,4,6-tribromophenol SUR		82	%			10 123		
		nitrobenzene-D5 SUR		62	%			35 114		
		2-fluorobiphenyl SUR		66	%			43 116		
		p-terphenyl-D14 SUR		77	%			33 141		

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit	
SW3005A6010C BLK5569		Silver		<	0.005	mg/L					
		Arsenic		<	0.008	mg/L					
		Barium		<	0.05	mg/L					
		Beryllium		<	0.004	mg/L					
		Cadmium		<	0.004	mg/L					
		Chromium		<	0.05	mg/L					
		Copper		<	0.05	mg/L					
		Nickel		<	0.05	mg/L					
		Lead		<	0.01	mg/L					
		Antimony		<	0.006	mg/L					
		Selenium		<	0.05	mg/L					
		Thallium		<	0.002	mg/L					
		Zinc		<	0.05	mg/L					
SW3005A6010C DUP5569		Silver	25147-001	<	0.005	mg/L			53	20	
		Arsenic	25147-001		0.022	mg/L			14	20	
		Cadmium	25147-001	<	0.004	mg/L			0	20	
		Chromium	25147-001	<	0.05	mg/L			108	20	
		Copper	25147-001	<	0.05	mg/L			3	20	
		Nickel	25147-001	<	0.05	mg/L			1	20	
		Lead	25147-001	<	0.01	mg/L			50	20	
		Zinc	25147-001	<	0.05	mg/L			1	20	
SW3005A6010C LCS5569		Silver		0.23	mg/L	0.25	93	80	120		
		Arsenic		0.49	mg/L	0.5	98	80	120		
		Barium		0.50	mg/L	0.5	100	80	120		
		Beryllium		0.49	mg/L	0.5	98	80	120		
		Cadmium		0.50	mg/L	0.5	100	80	120		
		Chromium		0.48	mg/L	0.5	97	80	120		
		Copper		0.51	mg/L	0.5	101	80	120		
		Nickel		0.50	mg/L	0.5	100	80	120		
		Lead		0.49	mg/L	0.5	99	80	120		
		Antimony		0.50	mg/L	0.5	99	80	120		
		Selenium		0.49	mg/L	0.5	99	80	120		
		Thallium		0.43	mg/L	0.5	86	80	120		
		Zinc		0.49	mg/L	0.5	98	80	120		
SW3005A6010C LCSD5569		Silver		0.24	mg/L	0.25	96	80	120	3	20
		Arsenic		0.50	mg/L	0.5	100	80	120	2	20
		Barium		0.50	mg/L	0.5	101	80	120	1	20
		Beryllium		0.50	mg/L	0.5	101	80	120	2	20
		Cadmium		0.51	mg/L	0.5	102	80	120	2	20
		Chromium		0.50	mg/L	0.5	99	80	120	3	20
		Copper		0.51	mg/L	0.5	103	80	120	2	20
		Nickel		0.51	mg/L	0.5	102	80	120	2	20
		Lead		0.50	mg/L	0.5	101	80	120	2	20
		Antimony		0.51	mg/L	0.5	102	80	120	2	20
		Selenium		0.51	mg/L	0.5	102	80	120	3	20
		Thallium		0.49	mg/L	0.5	98	80	120	13	20
		Zinc		0.50	mg/L	0.5	101	80	120	2	20

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3005A6010C MS5569	Silver		25147-001	0.31	mg/L	0.25	126	70 130		
	Arsenic		25147-001	0.69	mg/L	0.5	135	* 70 130		
	Cadmium		25147-001	0.62	mg/L	0.5	124	70 130		
	Chromium		25147-001	0.66	mg/L	0.5	132	* 70 130		
	Copper		25147-001	0.67	mg/L	0.5	132	* 70 130		
	Nickel		25147-001	0.64	mg/L	0.5	125	70 130		
	Lead		25147-001	0.63	mg/L	0.5	124	70 130		
	Zinc		25147-001	0.67	mg/L	0.5	131	* 70 130		
SW7470A	BLK5566	Mercury		< 0.0002	mg/L					
SW7470A	DUP5566	Mercury	25149-001	< 0.0002	mg/L					20
SW7470A	LCS5566	Mercury		0.0017	mg/L	0.002	86	80 120		
SW7470A	LCSD5566	Mercury		0.0017	mg/L	0.002	86	80 120	0	20
SW7470A	MS5566	Mercury	25149-001	0.0018	mg/L	0.002	90	75 125		

Absolute Resource

124 Heritage Avenue #168
Portsmouth, NH 03801
603-436-2001

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST**

25149

Project Name:
absoluteressourcesassociates.com

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S S O C I A T E S

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Sourceassociates.com

ANALYSIS REQUEST

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Absolute Resource a s s o c i a t e s										CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST		ANALYSIS REQUEST																																																																																																			
<p>Company Name: CREDERE ASSOCIATES</p> <p>Company Address: 776 MAIN ST, WESTBROOK, ME</p> <p>Report To: J. NEWCOMB/J. O'DONNELL</p> <p>Phone #: 207-828-1272</p> <p>Invoice To: T. PATTERN</p>										<p>Project Name: MERC</p> <p>Project #: 12001151</p> <p>Project Location: NH MA ME VT Other</p> <p>Protocol: RCRA SDWA NPDES NHDES OTHER</p> <p>Reporting QAPP GW-1 S-1 EPA DW Other</p> <p>Quote # _____ NH GREEODD Fund Pricing</p>		<p>124 Heritage Avenue #16 Portsmouth, NH 03801 603-436-2001 absoluteressourceassociates.com</p>																																																																																																			
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<p>TAT REQUESTED</p> <p>Priority (24 hr)* <input type="checkbox"/> Expedited (48 hr)* <input type="checkbox"/> Standard (10 Business Days) <input type="checkbox"/></p> <p>*Date Needed <u>10/17</u></p>										<p>SPECIAL INSTRUCTIONS</p> <p>See absoluteressourceassociates.com for sample acceptance policy and current accreditation lists.</p>																																																																																																					
<p>REPORTING INSTRUCTIONS</p> <p><input checked="" type="checkbox"/> PDF (e-mail address) <u>jnewcomb@credereassoc.com</u></p> <p><input type="checkbox"/> HARD COPY REQUIRED <input type="checkbox"/> FAX (FAX#) _____</p>										<p><input type="checkbox"/> OTHER (Specify)</p>																																																																																																					
<p>Relinquished by Sampler: <u>K. C.</u> Date <u>10/12/1515</u> Time <u>1515</u></p> <p>Received by: <u>J. O'Donnell</u></p>										<p>RECEIVED ON ICE <input type="checkbox"/> YES <input type="checkbox"/> NO</p> <p>TEMPERATURE <u>4</u> °C</p>																																																																																																					
<p>Relinquished by: _____ Date _____ Time _____</p>										<p>Received by Laboratory: _____ Date _____ Time _____</p>																																																																																																					
<p>OSD-01 Revision 08/23/12</p>										<p>25149</p>																																																																																																					

Credere-MERC samples group 1

CAPE Technologies EPA Method 4025m

Analysis completed October 10, 2012

Lab ID	Client ID	QA: Dup or	TEQ	Mean of	TEQ Sample Results	Corrected	Duplicate %CV	Notes
		pg/g spike	by 4025m	Dups		% spike recovery		
4-	CA-CC-4	0	44		30			unspiked sample
4-d	CA-CC-4	0 Dup	16	30			67	duplicate unspiked sample
4+	CA-CC-4	10	>45			xx		sample spiked with 10 pg/g 2378-TCDD
4+d	CA-CC-4	10 Dup	30	>38		0	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
5-	CA-CC-5	0	>45		>45			unspiked sample
5+	CA-CC-5	10	>45			xx		sample spiked with 10 pg/g 2378-TCDD
5+d	CA-CC-5	10 Dup	>45	>45		xx	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
6-	stack	0	11		12			unspiked sample
6-d	stack	0 Dup	13	12			11	duplicate unspiked sample
6+	stack	10	19			72		sample spiked with 10 pg/g 2378-TCDD
8-	CA-SS-2	0	>45		>200**			unspiked sample
8-d	CA-SS-2	0 Dup	>45	>45			xx	duplicate unspiked sample
8+	CA-SS-2	10	>45			xx		sample spiked with 10 pg/g 2378-TCDD
8+d	CA-SS-2	10 Dup	>45	>45		xx	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
11-	CA-SS-5	0	>45		>600**			unspiked sample
11+	CA-SS-5	10	>45			xx		sample spiked with 10 pg/g 2378-TCDD
11+d	CA-SS-5	10 Dup	>45	>45		xx	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
12-	CA-SS-DUP	0	>45		>400**			unspiked sample
12-d	CA-SS-DUP	0 Dup		xxx			xxx	duplicate unspiked sample
12+	CA-SS-DUP	10	>45			xx		sample spiked with 10 pg/g 2378-TCDD
MB-		0	2					unspiked method blank
MB-d		0 Dup	2	2			10	duplicate unspiked method blank
MB+		10	7			51		method blank spiked with 10 pg/g 2378-TCDD
MB+d		10 Dup	5	6		35	19	duplicate method blank spiked with 10 pg/g 2378-TCDD

Notes

All results are in pg/g (parts per trillion)

MB = Method Blank

d = Duplicate

"-" indicates unspiked sample results

"+" indicates matrix spike sample

"xx" indicates value which could not be calculated due to off scale high result

"xxx" indicates no valid result (sample lost during processing)

** reported result is based on two 5x serial dilutions of original sample residue

this procedure allows for rough estimation only

calibration adjustment factor includes correction for toluene residue effect as measured by evaporation controls

Credere-MERC samples group 2

CAPE Technologies EPA Method 4025m

Analysis completed October 10, 2012

Lab ID	Client ID	QA: Dup or	TEQ	Mean of	TEQ Sample Results	Corrected % spike recovery	Duplicate %CV	Notes
		pg/g spike	by 4025m	Dups				
1-	CA-SB-6	0	3		4			unspiked sample
1-d	CA-SB-6	0 Dup	4	4			14	duplicate unspiked sample
1+	CA-SB-6	10	11			71		sample spiked with 10 pg/g 2378-TCDD
2-	CA-SB-8 (0-2.5)	0	2		2			unspiked sample
2-d	CA-SB-8 (0-2.5)	0 Dup	2	2			10	duplicate unspiked sample
2+	CA-SB-8 (0-2.5)	10	7			53		sample spiked with 10 pg/g 2378-TCDD
2+d	CA-SB-8 (0-2.5)	10 Dup	6	6		43	10	duplicate sample spiked with 10 pg/g 2378-TCDD
3-	CA-SB-8 (5-10)	0	2		2			unspiked sample
3+	CA-SB-8 (5-10)	10	7			47		sample spiked with 10 pg/g 2378-TCDD
3+d	CA-SB-8 (5-10)	10 Dup	9	8		67	18	duplicate sample spiked with 10 pg/g 2378-TCDD
7-	CA-SS-1	0	>42		>42			unspiked sample
7+	CA-SS-1	10	>42			xx		sample spiked with 10 pg/g 2378-TCDD
7+d	CA-SS-1	10 Dup	>42	>42		xx	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
9-	CA-SS-3	0	10		13			unspiked sample
9-d	CA-SS-3	0 Dup	16	13			37	duplicate unspiked sample
9+	CA-SS-3	10	>42			xx		sample spiked with 10 pg/g 2378-TCDD
10-	CA-SS-4	0	34		>38			unspiked sample
10-d	CA-SS-4	0 Dup	>42	>38			xx	duplicate unspiked sample
10+	CA-SS-4	10	>42			xx		sample spiked with 10 pg/g 2378-TCDD
10+d	CA-SS-4	10 Dup	>42	>42		xx	xx	duplicate sample spiked with 10 pg/g 2378-TCDD
MB-		0	1					unspiked method blank
MB-d		0 Dup	2	2			2	duplicate unspiked method blank
MB+		10	6			46		method blank spiked with 10 pg/g 2378-TCDD
MB+d		10 Dup	8	7		68	22	duplicate method blank spiked with 10 pg/g 2378-TCDD

Notes

All results are in pg/g (parts per trillion)

MB = Method Blank

d = Duplicate

"-" indicates unspiked sample results

"+" indicates matrix spike sample

"xx" indicates value which could not be calculated due to off scale high result

calibration adjustment factor includes correction for toluene residue effect as measured by evaporation controls

Chain Of Custody Form



120 Thadeus Street
South Portland, ME 04106-6245
(207) 741-2995

Page 1 of 1

Project#: Proj. Name: MERC
Company: Credere Associates
Contact: Judd Newcomb
Address: 776 Main Street
Westbrook, ME 04092
Phone: 207-828-1272 PO# Quote #
Sampler (Signature):

Matrix Key:
C = Concrete
WP = Wipe
WW = Wastewater
SW = Surface Water
GW = Groundwater
DW = Drinking Water
S = Soil/Sludge
O = Oil
E = Extract
X = Other

Preservation

Samples were:

- 1) Shipped or hand-delivered on ice
- 2) Temp blank °C on ice
- 3) Received in good condition Y or N Y
- 4) pH checked by: NA
- 5) Labels checked by: NA

Container Key

P=plastic G=glass

Sample Identification	Sample Date	Sample Time	Analysis	Unpres	4°C	HNO ₃	H ₂ SO ₄	HCl	Methanol	Other	Matrix	Container number/type	pH	Cape Sample #
				X										
CA-SB-6	10/3/12	0850	Dioxin	X							S			1
CA-SB-8 (0-2.5")		0950									S			2
CA-SB-8 (5-10")		1000									S			3
CA-CC-4		1125									C			4
CA-CC-5		1148									C			5
Stack		1450									C			6
CA-SS-1		1120									S			7
CA-SS-2		1210									S			8
CA-SS-3		1215									S			9
CA-SS-4		1245									S			10
CA-SS-5		1300									S			11
CA-SS-Dup											S			12

Email Results to:
jnewcomb@crederelic.com

Comments / Instructions:

Ziplock bags

Turnaround Request

Standard <input checked="" type="checkbox"/>	Priority <input type="checkbox"/>
Due Date _____	Due Date _____
Lab Approval Required	

Project Requirements:

Relinquished By Sampler:

JBO

Received By:
R. Johnson

Date: 10/03/12 Time: 1610

Date: Time:

Relinquished By:
Relinquished By:
Relinquished By:

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

Judd Newcomb

CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: 12001150

Job ID: 25368

Date Received: 11/8/12

Project: MERC 12001150

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

A handwritten signature in black ink that appears to read "Sue Sylvester (for)".

Sue Sylvester
Principal, General Manager

Date of Approval: 11/13/2012
Total number of pages: 8

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SB-110 7.5-10	Solid	11/7/2012 15:00	25368-001	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-111 7.5-10	Solid	11/7/2012 15:05	25368-002	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-111 5-7.5	Solid	11/7/2012 15:10	25368-003	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SB-111 2.5-5	Solid	11/7/2012 15:15	25368-004	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G

Project ID: MERC 12001150

Job ID: 25368

Sample#: 25368-001

Sample ID: CA-SB-110 7.5-10

Matrix: Solid

Percent Dry: 85.3% Results expressed on a dry weight basis.

Sampled: 11/7/12 15:00		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	57	30-150	%	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A
decachlorobiphenyl SUR	50	30-150	%	1	JLZ	11/8/12	5625	11/9/12	14:23	SW3540C8082A

Sample#: 25368-002

Sample ID: CA-SB-111 7.5-10

Matrix: Solid

Percent Dry: 89.7% Results expressed on a dry weight basis.

Sampled: 11/7/12 15:05		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
PCB-1260	0.05	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	45	30-150	%	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A
decachlorobiphenyl SUR	66	30-150	%	1	JLZ	11/8/12	5625	11/9/12	14:53	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25368

Sample#: 25368-003

Sample ID: CA-SB-111 5-7.5

Matrix: Solid

Percent Dry: 87.2% Results expressed on a dry weight basis.

Sampled: 11/7/12 15:10		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1221	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1232	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1242	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1248	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1254	< 6.2	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
PCB-1260	55	6.2	ug/g	200	JLZ	11/8/12	5625	11/9/12	18:12	SW3540C8082A
Surrogate Recovery										
tetrachloro-m-xylene SUR										
DOR = Diluted out of range.										
DOR = Diluted out of range.										

Sample#: 25368-004

Sample ID: CA-SB-111 2.5-5

Matrix: Solid

Percent Dry: 93.6% Results expressed on a dry weight basis.

Sampled: 11/7/12 15:15		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1221	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1232	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1242	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1248	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1254	< 270	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
PCB-1260	6300	270	ug/g	10000	JLZ	11/8/12	5625	11/10/12	13:21	SW3540C8082A
Surrogate Recovery										
tetrachloro-m-xylene SUR										
DOR = Diluted out of range.										
DOR = Diluted out of range.										

Quality Control Report



124 Heritage Avenue Unit 16
Portsmouth, NH 03801
www.absoluteresourceassociates.com



Case Narrative
Lab # 25368

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 1 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

PCB: The surrogates were diluted out of the calibration range in the following samples: 25368-003 and -004.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3540C8082A	BLK5625	PCB-1016		<	0.02	ug/g				
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		<	0.02	ug/g				
		tetrachloro-m-xylene SUR		53	%			30 150		
		decachlorobiphenyl SUR		58	%			30 150		
SW3540C8082A	LCS5625	PCB-1016		0.15	ug/g	0.2	73	40 140		
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.14	ug/g	0.2	69	40 140		
		tetrachloro-m-xylene SUR		59	%			30 150		
		decachlorobiphenyl SUR		65	%			30 150		
SW3540C8082A	LCSD5625	PCB-1016		0.16	ug/g	0.2	81	40 140	10	30
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.16	ug/g	0.2	78	40 140	12	30
		tetrachloro-m-xylene SUR		60	%			30 150		
		decachlorobiphenyl SUR		69	%			30 150		

Absolute Resource

associates

124 Heritage Avenue #10
Portsmouth, NH 03801
603-436-2001
absoluteresourceassociates.com

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST****25368**

Company Name:

C R E E K A S S O C I A T E S

Company Address:

776 MAIN ST WESTBROOK ME

Report To:

JUDY NEWCOMBS

Phone #:

207-232-5387

Invoice To:

Same**ANALYSIS REQUEST**

Project Name: **MERC**
Project #: **10001150**
Project Location: NH MA ME VT Other
Protocol: RCRA SDWA NPDES
MCP NHDES OTHER
Reporting QAPP GW-1 S-1
Limits: EPA DW Other **75CA**

Quote #: **Lakeoke** NH GREE/DD
PO #: **10001150** Fund Pricing

Lab Sample ID	Field ID	# CONTAINERS	Matrix	Preservation Method	Sampling		Grab (G) or Composite (C)									
					WATER	SOLID	OTHER	HCl	HNO ₃	NaOH	MeOH	OTHER (Specify)	DATE	TIME	SAMPLER	
15368-1	CA-SB-111-25-1	1			X	X	X						4/1/12	1500	TR	
	CA-SB-111-25-1															
	CA-SB-111-25-1				X	X	X						1/5/12	1505		
	CA-SB-111-25-1				X	X	X						1/5/12	1510		
	CA-SB-111-25-1													1/5/12	1515	

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
 VOC 624 VOC BTEX MIBE, only VOC 8021VT
 VPH MADEP MEGRO GRO 8015
 VOC 524.2 VOC 524.2 NH List Gases-List:
 TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
 8270PAH 8270ABN 625 EDB 504.1
 8082 PCB 8081 Pesticides 608 Pest/PCB **3540 LCR**
 O&G 1664 Mineral O&G SM5520F
 pH BOD Conductivity Turbidity
 TSS TDS TS TVS Alkalinity
 RCRA Metals Priority Pollutant Metals TAL Metals
 Total Metals-list:
 Dissolved Metals-list:
 Ammonia COD TKN TN TON
 T-Phosphorus Phenols Bacteria P/A Bacteria MPN
 Cyanide Sulfide Nitrate + Nitrite Ortho P
 Nitrate Nitrite Chloride Sulfate Bromide Fluoride
 Corrosivity Reactive CN Reactive S- Ignitability/FP
 TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
Subcontract: TOC Grain Size TCLP Herbicides

TAT REQUESTED		SPECIAL INSTRUCTIONS	
Priority (24 hr)* <input type="checkbox"/>	Expedited (48 hr)* <input checked="" type="checkbox"/>	See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.	
REPORTING INSTRUCTIONS <input checked="" type="checkbox"/> PDF (e-mail address): JUNEWCOMBS@CREEK-RECORD.COM <input type="checkbox"/> OTHER (specify):			
REPORTING INSTRUCTIONS <input checked="" type="checkbox"/> PDF (e-mail address): JUNEWCOMBS@CREEK-RECORD.COM <input type="checkbox"/> OTHER (specify):		RECEIVED ON ICE <input checked="" type="checkbox"/> YES <input type="checkbox"/> NO	
RElinquished by Sampler: JUNEWCOMBS		Date 11/8/12	Time 1105
Relinquished by: JUNEWCOMBS		Date 11/8/12	Time 1105
Relinquished by: JUNEWCOMBS		Date 11/8/12	Time 1105

CUSTODY RECORD	
Relinquished by: JUNEWCOMBS	
Relinquished by: JUNEWCOMBS	
Relinquished by: JUNEWCOMBS	
OSD-01 Revision 12/23/10	

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

Judd Newcomb

CREDERE Associates
776 Main Street
Westbrook, ME 04092

PO Number: 12001150

Job ID: 25348

Date Received: 11/6/12

Project: MERC 12001150

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

A handwritten signature in black ink that appears to read "Sue Sylvester (for)".

Sue Sylvester
Principal, General Manager

Date of Approval: 11/14/2012
Total number of pages: 8

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Sample Association Table

Field ID	Matrix	Date-Time Sampled	Lab#	Analysis
CA-SS-6	Solid	11/5/2012 15:50	25348-001	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-7	Solid	11/5/2012 15:55	25348-002	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G
CA-SS-8	Solid	11/5/2012 16:00	25348-003	PCBs in soil by 8082 Percent Dry Matter for Sample Calc by SM2540B,G

Project ID: MERC 12001150

Job ID: 25348

Sample#: 25348-001

Sample ID: CA-SS-6

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 11/5/12 15:50		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	59	30-150	%	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A
decachlorobiphenyl SUR	55	30-150	%	1	JLZ	11/8/12	5625	11/9/12	11:31	SW3540C8082A

Sample#: 25348-002

Sample ID: CA-SS-7

Matrix: Solid

Percent Dry: 91.7% Results expressed on a dry weight basis.

Sampled: 11/5/12 15:55		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	66	30-150	%	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A
decachlorobiphenyl SUR	63	30-150	%	1	JLZ	11/8/12	5625	11/9/12	12:02	SW3540C8082A

Project ID: MERC 12001150

Job ID: 25348

Sample#: 25348-003

Sample ID: CA-SS-8

Matrix: Solid

Percent Dry: 93.3% Results expressed on a dry weight basis.

Sampled: 11/5/12 16:00		Reporting		Instr	Dil'n	Prep	Analysis			
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	Reference
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
Surrogate Recovery										
tetrachloro-m-xylene SUR	62	30-150	%	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A
decachlorobiphenyl SUR	59	30-150	%	1	JLZ	11/8/12	5625	11/9/12	12:32	SW3540C8082A

Quality Control Report



124 Heritage Avenue Unit 16
Portsmouth, NH 03801
www.absoluteresourceassociates.com



Case Narrative
Lab # 25348

Sample Receiving and Chain of Custody Discrepancies

Samples were received in acceptable condition, at 3 degrees C, on ice, and in accordance with sample handling, preservation and integrity guidelines.

Calibration

No exceptions noted.

Method Blank

No exceptions noted.

Surrogate Recoveries

No exceptions noted.

Laboratory Control Sample Results

No exceptions noted.

Matrix Spike/Matrix Spike Duplicate/Duplicate Results

Not requested for this project.

Other

Reporting Limits: Dilutions performed during the analysis are noted on the result pages.

No other exceptions noted.

- QC Report -

Method	QC ID	Parameter	Associated Sample	Result	Units	Amt Added	%R	Limits	RPD	RPD Limit
SW3540C8082A	BLK5625	PCB-1016		<	0.02	ug/g				
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		<	0.02	ug/g				
		tetrachloro-m-xylene SUR		53	%			30 150		
		decachlorobiphenyl SUR		58	%			30 150		
SW3540C8082A	LCS5625	PCB-1016		0.15	ug/g	0.2	73	40 140		
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.14	ug/g	0.2	69	40 140		
		tetrachloro-m-xylene SUR		59	%			30 150		
		decachlorobiphenyl SUR		65	%			30 150		
SW3540C8082A	LCSD5625	PCB-1016		0.16	ug/g	0.2	81	40 140	10	30
		PCB-1221		<	0.02	ug/g				
		PCB-1232		<	0.02	ug/g				
		PCB-1242		<	0.02	ug/g				
		PCB-1248		<	0.02	ug/g				
		PCB-1254		<	0.02	ug/g				
		PCB-1260		0.16	ug/g	0.2	78	40 140	12	30
		tetrachloro-m-xylene SUR		60	%			30 150		
		decachlorobiphenyl SUR		69	%			30 150		

Absolute Resource associates

124 Heritage Avenue #16
Portsmouth, NH 03801
603-436-2001

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

25348

PAGE 1 OF 1

absoluteresourceassociates.com

Company Name:

C R E O K E A S S O C I A T E S , L L C

Company Address:

776 MAIN ST WESTWICK NH

Report To:

TODD NEWCOMB

Phone #:

207 222-5382

Invoice To:

Same

ANALYSIS REQUEST

Project Name:	MERC
Project #:	1000150
Project Location:	NH MERC VT Other
Protocol:	RCRA SDWA NPDES
Reporting Limits:	MCP NHDES OTHER
Quote #:	CREOKE □ NH GREEODD Fund Pricing
PO #:	1000150

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
- VOC 624 VOC BTEX MTBE, only VOC 8021VT
- VPH MADEP MEGRO GRO 8015
- VOC 524.2 VOC 524.2 NH List Gases-List;
- TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
- 8270PAH 8270ABN 625 EDB 504.1
- 8082 PCB 8081 Pesticides 608 Pest/PCB **3540 Ext**
- O&G 1664 Mineral O&G SM5520F
- pH BOD Conductivity Turbidity
- TSS TDS TS TVS Alkalinity
- RCRA Metals Priority Pollutant Metals TAL Metals Hardness
- Total Metals-list:
- Dissolved Metals-list:
- Ammonia COD TKN TN TON TOC
- T-Phosphorus Phenols Bacteria P/A Bacteria MPN
- Cyanide Sulfide Nitrate + Nitrite Ortho P
- Nitrate Nitrite Chloride Sulfate Bromide Fluoride
- Corrosivity Reactive CN Reactive S- Ignitability/FP
- TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
- Subcontract: Grain Size Herbicides Formaldehyde

TAT REQUESTED (Lab Use Only)	Field ID	Matrix	Preservation Method	Sampling
Priority (24 hr)* <input type="checkbox"/>	CA-SS-6	WATER		
Expedited (48 hr)* <input type="checkbox"/>	CA-SS-2	SOLID		
Standard (10 Business Days) <input type="checkbox"/>	CA-SS-8	OTHER		

SPECIAL INSTRUCTIONS

See absoluteresourceassociates.com
for sample acceptance policy and
current accreditation lists.

REPORTING INSTRUCTIONS PDF (e-mail address)**TODDNEWCOMB@CREOKE.COM**

HARD COPY REQUIRED

FAX (FAX#)

OTHER (specify) **CREOKE LPP**

RECEIVED ON ICE YES NO
TEMPERATURE **13** °C

CUSTODY RECORD

Relinquished by: **TODD**

Date: **11/1/12**

Time: **11:05**

Received by: **TODD**

Date: **11/1/12**

Time: **11:05**

Received by Laboratory:

Date: **11/1/12**

Time: **11:05**

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

John Cressey
Summit Environmental Consultants
434 Cony Rd
Augusta, ME 04330

PO Number: 12-3259.1
Job ID: 25845
Date Received: 1/11/13

Project: Maine Energy 12-3259.1

Attached please find results for the analysis of the samples received on the date referenced above.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

A handwritten signature in black ink that appears to read "Sue Sylvester" followed by "(for)" in parentheses.

Sue Sylvester
Principal, General Manager

Date of Approval: 1/30/2013
Total number of pages: 12

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-001

Sample ID: SE-SB-201 (3"-4')

Matrix: Solid

Percent Dry: 92% Results expressed on a dry weight basis.

Sampled: 1/10/13 9:30		Reporting	Instr	Dil'n	Prep	Analysis			Reference	
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1254	1.7	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
PCB-1260	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	60	30-150	%	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082
decachlorobiphenyl SUR	61	30-150	%	5	JLZ	1/14/13	5777	1/16/13	21:08	SW3540C8082

Sample#: 25845-002

Sample ID: SE-SB-201 (4-5')

Matrix: Solid

Percent Dry: 86.1% Results expressed on a dry weight basis.

Sampled: 1/10/13 9:40		Reporting	Instr	Dil'n	Prep	Analysis			Reference	
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1254	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
PCB-1260	4.6	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	57	30-150	%	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082
decachlorobiphenyl SUR	58	30-150	%	5	JLZ	1/14/13	5777	1/16/13	21:39	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-003

Sample ID: SE-SB-202 (4"-2.5')

Matrix: Solid Percent Dry: 91.4% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:05		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1221	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1232	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1242	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1248	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1254	0.4	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
PCB-1260	< 0.1	0.1	ug/g	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	64	30-150	%	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082
decachlorobiphenyl SUR	66	30-150	%	1	JLZ	1/23/13	5802	1/24/13	15:52	SW3540C8082

Sample#: 25845-004

Sample ID: SE-SB-202 (4-5')

Matrix: Solid Percent Dry: 91.9% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:11		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1254	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
PCB-1260	< 0.2	0.2	ug/g	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	48	30-150	%	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082
decachlorobiphenyl SUR	29 *	30-150	%	5	JLZ	1/14/13	5777	1/16/13	22:10	SW3540C8082

* The surrogate showed recovery outside the acceptance limits. Reanalysis of the sample showed similar results. Matrix interference suspected.

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-005

Sample ID: SE-SB-203 (2"-2.5')

Matrix: Solid

Percent Dry: 94.4% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:25		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
PCB-1260	0.05	0.03	ug/g	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
Surrogate Recovery										
tetrachloro-m-xylene SUR										
	58	30-150	%	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082
decachlorobiphenyl SUR										
	66	30-150	%	1	JLZ	1/14/13	5777	1/15/13	18:21	SW3540C8082

Sample#: 25845-006

Sample ID: SE-SB-203 (3-4')

Matrix: Solid

Percent Dry: 84.3% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:31		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
PCB-1260	0.09	0.03	ug/g	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
Surrogate Recovery										
tetrachloro-m-xylene SUR										
	68	30-150	%	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082
decachlorobiphenyl SUR										
	57	30-150	%	1	JLZ	1/14/13	5777	1/16/13	20:38	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-007

Sample ID: SE-SB-203 (4-5')

Matrix: Solid

Percent Dry: 88.6% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:33		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1221	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1232	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1242	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1248	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1254	< 3.2	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
PCB-1260	85	3.2	ug/g	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	100	JLZ	1/14/13	5777	1/16/13	23:10	SW3540C8082

DOR = Diluted out of range.

Sample#: 25845-008

Sample ID: SE-SB-203 (5-7.5')

Matrix: Solid

Percent Dry: 87.3% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:37		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1221	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1232	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1242	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1248	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1254	< 0.6	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
PCB-1260	11	0.6	ug/g	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ	1/15/13	5777	1/18/13	13:27	SW3540C8082

DOR = Diluted out of range.

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-009

Sample ID: SE-SB-204 (3"-3')

Matrix: Solid

Percent Dry: 93.1% Results expressed on a dry weight basis.

Sampled: 1/10/13 10:56		Reporting	Instr	Dil'n	Prep	Analysis			Reference	
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	56	30-150	%	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082
decachlorobiphenyl SUR	51	30-150	%	1	JLZ	1/15/13	5777	1/17/13	1:42	SW3540C8082

Sample#: 25845-010

Sample ID: SE-SB-204 (3-5')

Matrix: Solid

Percent Dry: 88% Results expressed on a dry weight basis.

Sampled: 1/10/13 11:09		Reporting	Instr	Dil'n	Prep	Analysis			Reference	
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1221	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1232	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1242	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1248	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1254	< 0.3	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
PCB-1260	11	0.3	ug/g	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
Surrogate Recovery										
Limits										
tetrachloro-m-xylene SUR	DOR	30-150	%	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	10	JLZ	1/15/13	5777	1/18/13	12:57	SW3540C8082

DOR = Diluted out of range.

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-011

Sample ID: SE-SB-205 (3"-3')

Matrix: Solid

Percent Dry: 90.5% Results expressed on a dry weight basis.

Sampled: 1/10/13 11:24		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
PCB-1260	0.07	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	61	30-150	%	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082
decachlorobiphenyl SUR	46	30-150	%	1	JLZ	1/15/13	5777	1/17/13	2:43	SW3540C8082

Sample#: 25845-012

Sample ID: SE-SB-205 (3.5-7.5')

Matrix: Solid

Percent Dry: 94.9% Results expressed on a dry weight basis.

Sampled: 1/10/13 11:32		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
PCB-1260	0.30	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	60	30-150	%	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082
decachlorobiphenyl SUR	44	30-150	%	1	JLZ	1/15/13	5777	1/17/13	3:13	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-013

Sample ID: SE-SB-205 (10.5-12.5')

Matrix: Solid Percent Dry: 76.6% Results expressed on a dry weight basis.

Sampled: 1/10/13 11:47		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	53	30-150	%	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082
decachlorobiphenyl SUR	44	30-150	%	1	JLZ	1/15/13	5777	1/17/13	3:43	SW3540C8082

Sample#: 25845-014

Sample ID: SE-SB-210 (0-20")

Matrix: Solid Percent Dry: 83.3% Results expressed on a dry weight basis.

Sampled: 1/10/13 14:20		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1221	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1232	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1242	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1248	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1254	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
PCB-1260	< 0.04	0.04	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	51	30-150	%	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082
decachlorobiphenyl SUR	34	30-150	%	1	JLZ	1/15/13	5777	1/17/13	4:14	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25845

Sample#: 25845-015

Sample ID: SE-SB-211 (2"-3.5')

Matrix: Solid Percent Dry: 89.4% Results expressed on a dry weight basis.

Sampled: 1/10/13 14:58		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
PCB-1260	0.23	0.0	ug/g	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
Surrogate Recovery										
tetrachloro-m-xylene SUR										
	63	30-150	%	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082
decachlorobiphenyl SUR										
	33	30-150	%	1	JLZ	1/15/13	5777	1/17/13	4:44	SW3540C8082

Sample#: 25845-016

Sample ID: SE-SB-211 (4-9')

Matrix: Solid Percent Dry: 87.2% Results expressed on a dry weight basis.

Sampled: 1/10/13 15:03		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1221	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1232	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1242	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1248	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1254	< 1.6	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
PCB-1260	48	1.6	ug/g	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
Surrogate Recovery										
tetrachloro-m-xylene SUR										
	DOR	30-150	%	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082
decachlorobiphenyl SUR										
	DOR	30-150	%	50	JLZ	1/15/13	5777	1/18/13	13:58	SW3540C8082

DOR = Diluted out of range.

Absolute Resource associates

124 Heritage Avenue #16
Portsmouth, NH 3801
603-436-2001
absoluteresourceassociates.com

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST 25845

PAGE / OF 3

Company Name:
SUMMIT ENVIRONMENTAL
Company Address:
640 MAIN ST. LEXINGTON, ME

Report To:
John Cresssey
Phone #:
(207) 745-6659

Invoice To:
SAME

Project Name: **MAINE ENERGY**

Project #: **12-3259.1**

Project Location: NH M **ME** VT Other

Protocol: **RHCA MCP** SDWA NPDES OTHER

Reporting Limits: QAPP GW-1 S-1 EPA DW Other

PO # **12-3259.1**
Quote # **16434** □ NH GREE/ODD Fund Pricing

ANALYSIS REQUEST

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
- VOC 624 VOC BTEX MTBE, only VOC 8021VT
- VPH MADEP MEGRO GRO 8016
- VOC 524.2 VOC 524.2 NH List Gases-List
- TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
- 8270PAH 8270ABN 625 EDB 504.1
- 82802 PCB 8081 Pesticides 608 Pest/PCB
- O&G 1664 Mineral O&G SM5520F
- pH BOD Conductivity Turbidity
- TSS TDS TS TVS Alkalinity
- RCRA Metals Priority Pollutant Metals TAL Metals Hardness
- Total Metals-list:
- Dissolved Metals-list:
- Ammonia COD TKN TN TON TOC
- T-Phosphorus Phenols Bacteria P/A Bacteria MPN
- Cyanide Sulfide Nitrate + Nitrite Ortho P
- Nitrate Nitrite Chloride Sulfate Bromide Fluoride
- Corrosivity Reactive CN Reactive S- Ignitability/FP
- TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
- Subcontract: Grain Size Herbicides Formaldehyde

Lab Sample ID

Field ID

Matrix

Preservation Method

Sampling

Lab Sample ID (Lab Use Only)	Field ID # CONTAINERS	Matrix	Preservation Method	Sampling	6°C		SAMPLER								
					WATER	SOLID		OTHER	HCl	HNO ₃	H ₂ SO ₄	NaOH	MeOH	OTHER (Specify)	DATE
25615-01	SE-SB-201 (3x1)	1	X	X					X	1-10	9:30	JRC			G
-02	SE-SB-201 (4-5)	1	X	X					X	1-10	9:40	JRC			G
-03	SE-SB-202 (4-5)	1	X	X					X	1-10	10:05	JRC			G
-04	SE-SB-202 (4-5)	1	X	X					X	1-10	10:11	JRC			G
-05	SE-SB-203 (3-25)	1	X	X					X	1-10	10:25	JRC			G
-06	SE-SB-203 (3-4)	1	X	X					X	1-10	10:31	JRC			G
-07	SE-SB-203 (4-5)	1	X	X					X	1-10	10:33	JRC			G
-08	SE-SB-203 (5-7)	1	X	X					X	1-10	10:37	JRC			G
-09	SE-SB-204 (3-3)	1	X	X					X	1-10	10:56	JRC			G
-10	SE-SB-204 (3-5)	1	X	X					X	1-10	11:09	JRC			G
-11	SE-SB-205 (4-3)	1	X	X					X	1-10	11:24	JRC			G

TAT REQUESTED

Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)
*Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

REPORTING INSTRUCTIONS

PDF (e-mail address)

JRCESSEY@COMCAST.NET

HARD COPY REQUIRED

FAX (FAX#)

OTHER (specify)

SPECIAL INSTRUCTIONS

SOXHLET EXTRACTION

CUSTODY RECORD

Relinquished by: *JRC*

Date: *1-11-13*

Time: *12:45*

Received by: *JRC*

Date: *1-11-13*

Time: *12:45*

Received by Laborator: _____

Date: _____

Time: _____

Absolute Resource

124 Heritage Avenue #16
Portsmouth, NH 03801
603-436-2001
absoluteresourceassociates.com

ANALYSIS REQUEST

CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST
25845

PAGE 2 OF 3

Company Name:
SUMMIT ENVIRONMENTAL
Company Address:
640 Main St. Lewiston, ME

Report To:
Jesse Cresser

Phone #:
(207)745-6001

Invoice To:
Jesse Cresser

Lab Sample ID	Field ID	Matrix	Preservation Method	Sampling
385-12	SE-SB-205(3.5#)	1	# CONTAINERS	
-13	SE-SB-205(6.5#)	1	WATER	X
-14	SE-SB-210(6.25#)	1	SOLID	X
-15	SE-SB-210(2.5#)	1	OTHER	X
-16	SE-SB-210(4.9#)	1	HCl	X
-17	SE-D-01	1	HNO ₃	X
-18	SE-D-02	1	H ₂ SO ₄	X
-19	SE-D-03	1	NaOH	X
-20	SE-D-04	1	MeOH	X
-21	SE-D-05	1	OTHER (Specify)	X
-22	SE-D-06	1	DATE	1-10
			TIME	1130 JMC
			SAMPLER	JMC

Protocol: <input checked="" type="checkbox"/> RCRA <input type="checkbox"/> SDWA <input type="checkbox"/> NPDES <input type="checkbox"/> MDEP <input type="checkbox"/> NHDES <input type="checkbox"/> OTHER
Reporting Limits: <input type="checkbox"/> QAPP <input type="checkbox"/> EPA DW <input type="checkbox"/> Other
Quote #: 6434 <input type="checkbox"/> NH GREE/ODD Fund Pricing PO #: 12-325911

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
- VOC 624 VOC BTEX MIBE, only VOC 8021VT
- VPH MADEP MEGRO GRO 8015
- VOC 524.2 VOC 524.2 NH List Gases-List:
- TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
- 8270PAH 8270ABN 625 EDB 504.1
- 8082 PCB 8081 Pesticides 608 Pest/PCB
- O&G 1664 Mineral O&G SM5520F
- pH BOD Conductivity Turbidity
- TSS TDS TS TVS Alkalinity
- RCRA Metals Priority Pollutant Metals TAL Metals Hardness
- Total Metals-List:
- Dissolved Metals-List:
- Ammonia COD TKN TN TON TOC
- T-Phosphorus Phenols Bacteria P/A Bacteria MPN
- Cyanide Sulfide Nitrate + Nitrite Ortho P
- Nitrate Nitrite Chloride Sulfate Bromide Fluoride
- Corrosivity Reactive CN Reactive S- Ignitability/FP
- TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
- Subcontract: Grain Size Herbicides Formaldehyde

DIOXIN TEQ

Grab (G) or Composite (C)

*TAT REQUESTED
Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)
Date Needed _____

REPORTING INSTRUCTIONS
See absoluteresourceassociates.com
for sample acceptance policy and
current accreditation lists.

HARD COPY REQUIRED FAX (FAX#)
Relinquished by: *Jesse Cresser* Date: **1-11-13** Time: **1224** Received by: *Jesse Cresser*

OTHER (specify)
RECEIVED ON ICE YES NO
TEMPERATURE **3** °C

CUSTODY
RECORD

Relinquished by: _____ Date: _____ Time: _____ Received by Laboratory: _____

OSD-01 Revision 08/23/12

Absolute Resource

associates

124 Heritage Avenue #16
Portsmouth, NH 03801
603-436-2001

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

PAGE 3 OF 3

Company Name:
SUMMIT ENVIRONMENTAL
Company Address:
640 Main St Lewiston ME 04240
Report To:
John Cresser
Phone #:
(302) 745-6609
Invoice To:
JAME

Project Name: **MAINE ENERGY**
Project #: _____

Project Location: **NH MA ME VT**
Protocol: **RCCA MCP NHDES OTHER**
Reporting Limits: **QAPP GW-1 S-1 EPA DW Other**
Quote #: _____
PO #: _____

Lab Sample ID (Lab Use Only)	Field ID	# CONTAINERS	Matrix	Preservation Method	Sampling					
					WATER	SOLID	OTHER	DATE	TIME	SAMPLER
35385-23	SE-D-07	1	X	HCl		X		1-11	9:47 AM	SCD
-24	SE-D-08	1	X	HNO ₃		X		1-11	10:00	
-25	SE-D-09	1	X	H ₂ SO ₄		X		1-11	9:55	
-26	SE-D-10	1	X	NaOH		X		1-11	10:10	
-27	SE-D-11	1	X	MeOH		X		1-11	10:32	
-28	SE-D-12	1	X	OTHER (Specify)		X		1-11	10:40	
-29	SE-D-13	1	X			X		1-11	10:49	
-30	SE-D-14	1	X			X		1-11	10:56	
-31	SE-D-15	1	X			X		1-11	11:10	

- VOC 8250 VOC 8260 NHDES VOC 8260 MADEP
 - VOC 624 VOC BTEX MIBE, only VOC 8021VT
 - VPH MADEP MEGRO GRO 8015
 - VOC 524.2 VOC 524.2 NH List Gases-List:
 - TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
 - 8270PAH 8270ABN 625 EDB 504.1
 - 8082 PCB 8081 Pesticides 608 Pest/PCB
 - O&G 1664 Mineral O&G SM5520F
 - pH BOD Conductivity Turbidity
 - TSS TDS TS TVS Alkalinity
 - RCRA Metals Priority Pollutant Metals TAL Metals Hardness
 - Total Metals-list:
 - Dissolved Metals-list:
 - Ammonia COD TKN TN TON TOC
 - T-Phosphorus Phenols Bacteria P/A Bacteria MPN
 - Cyanide Sulfide Nitrate + Nitrite Ortho P
 - Nitrate Nitrite Chloride Sulfate Bromide Fluoride
 - Corrosivity Reactive CN Reactive S- Ignitability/FP
 - TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
 - Subcontract: Grain Size Herbicides Formaldehyde
- Grab (G) or Composite (C)
- DIOXIN TEQ

TAT REQUESTED

Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)

*Date Needed _____

See absoluteresourceassociates.com for sample acceptance policy and current accreditation lists.

SPECIAL INSTRUCTIONS

REPORTING INSTRUCTIONS PDF (e-mail address) _____

HARD COPY REQUIRED FAX (FAX#) _____

RECEIVED ON ICE YES NO
TEMPERATURE 3 °C

Date 1-11-13 Time 10:24

Received by: J. Cresser Date Time

Received by: _____ Date Time

Relinquished by: _____

Date Time

Reinquished by: _____

Date Time

Received by Laboratory: _____

Date Time

QSD-01 Revision 08/23/12

Laboratory Report



Absolute Resource associates

124 Heritage Avenue Portsmouth NH 03801

John Cressey
Summit Environmental Consultants
434 Cony Rd
Augusta, ME 04330

PO Number: 12-3259.1
Job ID: 25852
Date Received: 1/11/13

Project: Maine Energy 12-3259.1

Attached please find results for the analysis of the samples received on the date referenced above.
The following report has been re-issued to correct a sample ID, as requested by the customer.

Unless otherwise noted in the attached report, the analyses performed met the requirements of Absolute Resource Associates' Quality Assurance Plan. The Standard Operating Procedures are based upon USEPA SW-846, USEPA Methods for Chemical Analysis of Water and Wastewater, Standard Methods for the Examination of Water and Wastewater and other recognized methodologies. The results contained in this report pertain only to the samples as indicated on the chain of custody.

Absolute Resource Associates maintains certification with the agencies listed below.

We appreciate the opportunity to provide laboratory services. If you have any questions regarding the enclosed report, please contact the laboratory and we will be glad to assist you.

Sincerely,
Absolute Resource Associates

A handwritten signature in black ink that appears to read "Sue Sylvester (for)".

Sue Sylvester
Principal, General Manager

Date of Approval: 2/15/2013
Total number of pages: 10

Absolute Resource Associates Certifications

New Hampshire 1732
Maine NH903

Massachusetts M-NH902

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-001

Sample ID: SE-SB-206 (3"-2')

Matrix: Solid

Percent Dry: 91.8% Results expressed on a dry weight basis.

Sampled: 1/10/13 11:58		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1254	< 0.2	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
PCB-1260	1.7	0.2	ug/g	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	51	30-150	%	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082
decachlorobiphenyl SUR	54	30-150	%	5	JLZ	1/15/13	5777	1/18/13	12:26	SW3540C8082

Sample#: 25852-002

Sample ID: SE-SB-206 (2'-12')

Matrix: Solid

Percent Dry: 93% Results expressed on a dry weight basis.

Sampled: 1/10/13 12:01		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
PCB-1260	0.31	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	61	30-150	%	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082
decachlorobiphenyl SUR	76	30-150	%	1	JLZ	1/16/13	5786	1/18/13	15:36	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-003

Sample ID: SE-SB-207 (3"-3')

Matrix: Solid

Percent Dry: 92.5% Results expressed on a dry weight basis.

Sampled: 1/10/13 12:48		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	57	30-150	%	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082
decachlorobiphenyl SUR	62	30-150	%	1	JLZ	1/16/13	5786	1/18/13	15:06	SW3540C8082

Sample#: 25852-004

Sample ID: SE-SB-207 (3'-4.5')

Matrix: Solid

Percent Dry: 90.3% Results expressed on a dry weight basis.

Sampled: 1/10/13 12:51		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1221	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1232	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1242	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1248	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1254	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
PCB-1260	6.3	0.4	ug/g	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	10	JLZ	1/16/13	5786	1/18/13	23:23	SW3540C8082

DOR = Diluted out of range.

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-005

Sample ID: SE-SB-207 (5-6')

Matrix: Solid

Percent Dry: 88.5% Results expressed on a dry weight basis.

Sampled: 1/10/13 12:57

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
PCB-1016	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1254	< 0.2	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
PCB-1260	2.3	0.2	ug/g	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
Surrogate Recovery									
tetrachloro-m-xylene SUR	46	30-150	%	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082
decachlorobiphenyl SUR	40	30-150	%	5	JLZ 1/16/13	5786	1/18/13	23:53	SW3540C8082

Sample#: 25852-006

Sample ID: SE-SB-208 (3"-4')

Matrix: Solid

Percent Dry: 93.9% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:10

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Analysis			Reference
	Result	Limit				Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
PCB-1260	< 0.03	0.03	ug/g	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
Surrogate Recovery									
tetrachloro-m-xylene SUR	60	30-150	%	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082
decachlorobiphenyl SUR	66	30-150	%	1	JLZ 1/16/13	5786	1/18/13	16:47	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-007

Sample ID: SE-SB-208 (4-6')

Matrix: Solid

Percent Dry: 86.7% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:12		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1221	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1232	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1242	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1248	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1254	< 0.2	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
PCB-1260	2.5	0.2	ug/g	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	82	30-150	%	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082
decachlorobiphenyl SUR	54	30-150	%	5	JLZ	1/16/13	5786	1/21/13	16:10	SW3540C8082

Sample#: 25852-008

Sample ID: SE-SB-208 (6-8')

Matrix: Solid

Percent Dry: 84.5% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:16		Reporting Limit		Instr Dil'n		Prep Date	Analysis			Reference
Parameter	Result	Units	Factor	Analyst	Batch	Date	Time			
PCB-1016	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1221	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1232	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1242	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1248	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1254	< 0.4	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
PCB-1260	5.0	0.4	ug/g	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	10	JLZ	1/16/13	5786	1/21/13	16:40	SW3540C8082

DOR = Diluted out of range.

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-009

Sample ID: SE-SB-208 (8-10')

Matrix: Solid

Percent Dry: 77.6% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:23		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1221	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1232	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1242	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1248	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1254	< 2.1	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
PCB-1260	25	2.1	ug/g	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	50	JLZ	1/17/13	5786	1/21/13	17:41	SW3540C8082

DOR = Diluted out of range.

Sample#: 25852-010

Sample ID: SE-SB-209 (2"-3')

Matrix: Solid

Percent Dry: 92.4% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:31		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
PCB-1260	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	63	30-150	%	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082
decachlorobiphenyl SUR	70	30-150	%	1	JLZ	1/17/13	5786	1/18/13	21:52	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-011

Sample ID: SE-SB-209 (3-9')

Matrix: Solid

Percent Dry: 87.1% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:37		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1221	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1232	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1242	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1248	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1254	< 360	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
PCB-1260	3500	360	ug/g	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	DOR	30-150	%	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082
decachlorobiphenyl SUR	DOR	30-150	%	10000	JLZ	1/17/13	5786	1/21/13	15:36	SW3540C8082

DOR = Diluted out of range.

Sample#: 25852-012

Sample ID: SE-SB-209 (9-12.5')

Matrix: Solid

Percent Dry: 89.1% Results expressed on a dry weight basis.

Sampled: 1/10/13 13:45		Reporting		Instr	Dil'n	Prep	Analysis			Reference
Parameter	Result	Limit	Units	Factor	Analyst	Date	Batch	Date	Time	
PCB-1016	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1221	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1232	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1242	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1248	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1254	< 0.03	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
PCB-1260	0.13	0.03	ug/g	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
Surrogate Recovery		Limits								
tetrachloro-m-xylene SUR	74	30-150	%	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082
decachlorobiphenyl SUR	82	30-150	%	1	JLZ	1/17/13	5786	1/21/13	13:21	SW3540C8082

Project ID: Maine Energy 12-3259.1

Job ID: 25852

Sample#: 25852-013

Sample ID: SE-SB-212 (2-20")

Matrix: Solid

Percent Dry: 91.1% Results expressed on a dry weight basis.

Sampled: 1/10/13 14:45

Parameter	Reporting		Instr Dil'n	Analyst	Prep Date	Batch	Analysis			Reference
	Result	Limit					Date	Time		
PCB-1016	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1221	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1232	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1242	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1248	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1254	< 0.6	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
PCB-1260	7.1	0.6	ug/g	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
Surrogate Recovery										
tetrachloro-m-xylene SUR	DOR	30-150	%	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	
decachlorobiphenyl SUR	DOR	30-150	%	20	JLZ 1/17/13	5786	1/21/13	17:11	SW3540C8082	

DOR = *Diluted out of range.*

SET #2 - ANALYZE ONLY AFTER TALKING WITH SUMMIT

PAGE 1 OF 2

Absolute Resource associates

124 Heritage Avenue #16
Portsmouth, NH 03801
603-436-2001
absoluteresourceassociates.com

CHAIN-OF-CUSTODY RECORD AND ANALYSIS REQUEST

25852

ANALYSIS REQUEST

Company Name: SUMMIT ENVIRONMENTAL	Project Name: MAINE ENERGY
Company Address: 140 MAIN ST. LEXINGTON, ME 04441	Project #: 12-325911
Report To: John Cresser	Project Location: NH MAINE
Phone #: (207) 745-6009	Protocol: <input checked="" type="checkbox"/> RCRA <input type="checkbox"/> SDWA <input type="checkbox"/> NPDES <input type="checkbox"/> MCP <input type="checkbox"/> NHDES <input type="checkbox"/> OTHER
Invoice To: SUMMIT	Reporting Limits: <input type="checkbox"/> QAPP <input type="checkbox"/> GW-1 <input type="checkbox"/> S-1 <input type="checkbox"/> EPA DW <input type="checkbox"/> Other
	Quote #: 60434 <input type="checkbox"/> NH GREE/ODD Fund Pricing PO #: 12-325911

Lab Sample ID	Field ID	Matrix	Preservation Method	Sampling
(Lab Use Only)		# CONTAINERS		
233301	SE-SB-206(3-1)	1	WATER	
.02	SE-SB-206(3-1)	1	SOLID	
.03	SE-SB-207(3-3)	1	OTHER	
.04	SE-SB-207(3-4)	1	HCl	
.05	SE-SB-207(3-6)	1	HNO ₃	
.06	SE-SB-208(3-4)	1	H ₂ SO ₄	
.07	SE-SB-208(4-6)	1	NaOH	
.08	SE-SB-208(6-3)	1	MeOH	
.09	SE-SB-208(8-10)	1	OTHER (Specify)	
.10	SE-SB-208(9-3)	1		
.11	SE-SB-208(3-9)	1		

- VOC 8260 VOC 8260 NHDES VOC 8260 MADEP
- VOC 824 VOC BTEX MIBE, only VOC 8021VT
- VPH MADEP MEGRO GRO 8015
- VOC 524.2 VOC 524.2 NH List Gases-List:
- TPH DRO 8015 MEDRO EPH MADEP TPH Fingerprint
- 8270PAH 8270ABN 625 EDB 504.1
- 8082 PCB 8081 Pesticides 608 Pest/PCB
- O&G 1664 Mineral O&G SM5520F
- pH BOD Conductivity Turbidity
- TSS TDS TS TVS Alkalinity
- RCRA Metals Priority Pollutant Metals TAL Metals Hardness
- Total Metals-list:
- Dissolved Metals-list:
- Ammonia COD TKN TN TON TOC
- T-Phosphorus Phenols Bacteria P/A Bacteria MPN
- Cyanide Sulfide Nitrate + Nitrite Ortho P
- Nitrate Nitrite Chloride Sulfate Bromide Fluoride
- Corrosivity Reactive CN Reactive S- Ignitability/FP
- TCLP Metals TCLP VOC TCLP SVOC TCLP Pesticide
- Subcontract: Grain Size Herbicides Formaldehyde

per case from John Wiles,
proceed w/ analysis
of these samples

RECEIVED ON ICE YES NO
TEMPERATURE **3** °C
Grab (G) or Composite (C)

TAT REQUESTED

Priority (24 hr)*
Expedited (48 hr)*
Standard (10 Business Days)

REPORTING INSTRUCTIONS HARD COPY REQUIRED FAX (FAX#) _____

See absoluteresourceassociates.com
for sample acceptance policy and
current accreditation lists.

ANALYZE ONLY AFTER AUTHORIZATION FROM JOHN WILES

John Wiles
per case from John Wiles,
proceed w/ analysis
of these samples

*Date Needed _____

CUSTODY RECORD

Relinquished by: *J. Cresser* Date: **1-11-13** Time: **1224**
Received by: _____

Relinquished by: *J. Cresser* Date: _____ Time: _____
Received by: _____

OSD-01 Revision 08/23/12

LESS

PAGE 2 OF 2

Absolute Resource

[absoluteresourceassociates.com](http://www.absoluteresourceassociates.com)

Sumit Environmental
Company Limited

Company Address:
640 MAIN ST. LEXINGTON, ME

Report To:

Diana M.
DHS Lesser

Phone #:

Invoice To:

JOURNAL OF POLYMER SCIENCE: PART A

TABLE III
CONTINUATION

(Lab Use Only) DI DI # CONC WATER SOLID OTHER HCl ING.

2880-012 SE-SB-20(F-10.3) 1 X

13 SE-SS-212(2-25) 1 X

ANSWER

1000

ANSWER

ANSWER

ANSWER

卷之三

TAT REQUESTED Priority (24 hr)* See absoluteresourceassociates.com for sample acceptance policy and

Expedited (48 hr)* **Standard** **current accreditation lists.**

REPORTING INSTRUCTION

DISMANTLED Relinquished by Sample:

RECORDED BY _____
Relinquished by: _____

RECORD
OSD-01 Revision 08/23/12
Relinquished by: *[Signature]*

**CHAIN-OF-CUSTODY RECORD
AND ANALYSIS REQUEST**

25852

February 27, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74873
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples
PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
Chromatograms
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:

No exceptions.

PCBs by EPA Method 8082:

No results were reported below the quantitation limit.

All samples except 74873-1, 74873-3 and 74873-19 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument.

Sample 74873-18 had high recoveries for surrogate Tetrachloro-m-xylene. Decachlorobiphenyl surrogate recoveries were in control. Results were reported with a comment to this affect. Some samples had interference that prevented the determination of a surrogate recovery. Results were reported off the column without interferences when possible, or the sample was reported with a comment to this affect.

The MS/MSD analyzed on sample 74873-1 had some high RPDs (see form 3). The laboratory control samples (L022013PSOX2, RR/LD022013PSOX2, RR) were in control. Results were reported without qualification.

The MS/MSD analyzed on sample 74873-16 was only assessed for precision of PCB 1260 due to the parent sample having concentrations of PCB 1260 that exceeded the calibration range of the instrument. The RPDs for the MS/MSD did not meet acceptance criteria. The laboratory control samples (L022513PSOX/LD022513PSOX) were in control for all recoveries and RPDs. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC



Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
Summit Environmental Consultants Inc.
640 Main Street
Lewiston ME 04240

Report Number: 74873

Revision: Rev. 0

Re: MAINE ENERGY (Project No: 12-3259.1)

Enclosed are the results of the analyses on your sample(s). Samples were received on 20 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Sample Analysis: The attached pages detail the Client Sample IDs, Lab Sample IDs, and Analyses requested

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

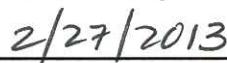
If you have any questions on these results, please do not hesitate to contact us.

Authorized signature



Stephen L. Knollmeyer Lab. Director

Date



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CLIENT: Summit Environmental
 Consultants Inc.

REPORT NUMBER: 74873

REV: Rev. 0

PROJECT: MAINE ENERGY (Project No: 12-3259.1)

<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74873-1	02/19/13	SE-SB-216 (2.5"-24")	EPA 8082 (PCBs only)	
74873-2	02/19/13	SE-SB-216 (2-4')	EPA 8082 (PCBs only)	
74873-3	02/19/13	SE-SB-216 (4-6')	EPA 8082 (PCBs only)	
74873-4	02/19/13	SE-SB-215 (0-2')	EPA 8082 (PCBs only)	
74873-5	02/19/13	SE-SB-215 (2-4')	EPA 8082 (PCBs only)	
74873-6	02/19/13	SE-SB-215 (4-6')	EPA 8082 (PCBs only)	
74873-7	02/19/13	SE-SB-215 (6-8")	EPA 8082 (PCBs only)	
74873-8	02/19/13	SE-SB-215 (8-10')	EPA 8082 (PCBs only)	
74873-9	02/19/13	SE-SB-214 (0-2')	EPA 8082 (PCBs only)	
74873-10	02/19/13	SE-SB-214 (2-4')	EPA 8082 (PCBs only)	
74873-11	02/19/13	SE-SB-214 (4-6')	EPA 8082 (PCBs only)	
74873-12	02/19/13	SE-SB-214 (6-7")	EPA 8082 (PCBs only)	
74873-13	02/19/13	SE-SB-213 (0-2')	EPA 8082 (PCBs only)	
74873-14	02/19/13	SE-SB-213 (2-4')	EPA 8082 (PCBs only)	
74873-15	02/19/13	SE-SB-213 (4-6')	EPA 8082 (PCBs only)	
74873-16	02/19/13	SE-SB-213 (6-7.5')	EPA 8082 (PCBs only)	
74873-17	02/19/13	SE-SB-217 (0-2')	EPA 8082 (PCBs only)	
74873-18	02/19/13	SE-SB-217 (2-4')	EPA 8082 (PCBs only)	
74873-19	02/19/13	SE-SB-217 (4-6')	EPA 8082 (PCBs only)	
74873-20	02/19/13	SE-SB-217 (8-10')	EPA 8082 (PCBs only)	

Surrogate Compound Limits

	Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water				
1,4-Difluorobenzene		70-130		EPA 524.2
Bromofluorobenzene		70-130		
1,2-Dichlorobenzene-d4		70-130		
Volatile Organic Compounds				
1,2-Dichloroethane-d4		70-120	70-120	EPA 624/8260B
Toluene-d8		85-120	85-120	
Bromofluorobenzene		75-120	75-120	
Semi-Volatile Organic Compounds				
2-Fluorophenol		20-110	35-105	EPA 625/8270C
d5-Phenol		15-110	40-100	
d5-nitrobenzene		40-110	35-100	
2-Fluorobiphenyl		50-110	45-105	
2,4,6-Tribromophenol		40-110	40-125	
d14-p-terphenyl		50-130	30-125	
PAH's by SIM				
d5-nitrobenzene		21-110	35-110	EPA 8270C
2-Fluorobiphenyl		36-121	45-105	
d14-p-terphenyl		33-141	30-125	
Pesticides and PCBs				
2,4,5,6-Tetrachloro-m-xylene (TCX)		46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)		40-135	40-130	
Herbicides				
Dichloroacetic acid (DCAA)		30-150	30-150	
Gasoline Range Organics/TPH Gasoline				
Trifluorotoluene TFT (FID)		60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)		60-140	60-140	
Trifluorotoluene TFT (PID)		60-140	60-140	
Bromofluorobenzene (BFB) (PID)		60-140	60-140	
Diesel Range Organics/TPH Diesel				
m-terphenyl		60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons				
2,5-Dibromotoluene (PID)		70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)		70-130	70-130	
Extricable Petroleum Hydrocarbons				
1-chloro-octadecane (aliphatic)		40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)		40-140	40-140	
2-Fluorobiphenyl (Fractionation)		40-140	40-140	
2-Bromonaphthalene (fractionation)		40-140	40-140	

PCB
DATA SUMMARIES

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-216 (2.5"-24")

Lab Sample ID:	74873-1 RR
Matrix:	Solid
Percent Solid:	95
Dilution Factor:	1.0
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/20/13
Analysis Date:	02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	65

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	90	%
Decachlorobiphenyl	75	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-1,RR,,A/C

Column ID: 0.25 mm

Data File: M67211.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	65	64	2.9	

Column to be used to flag RPD values greater than QC limit of 40%

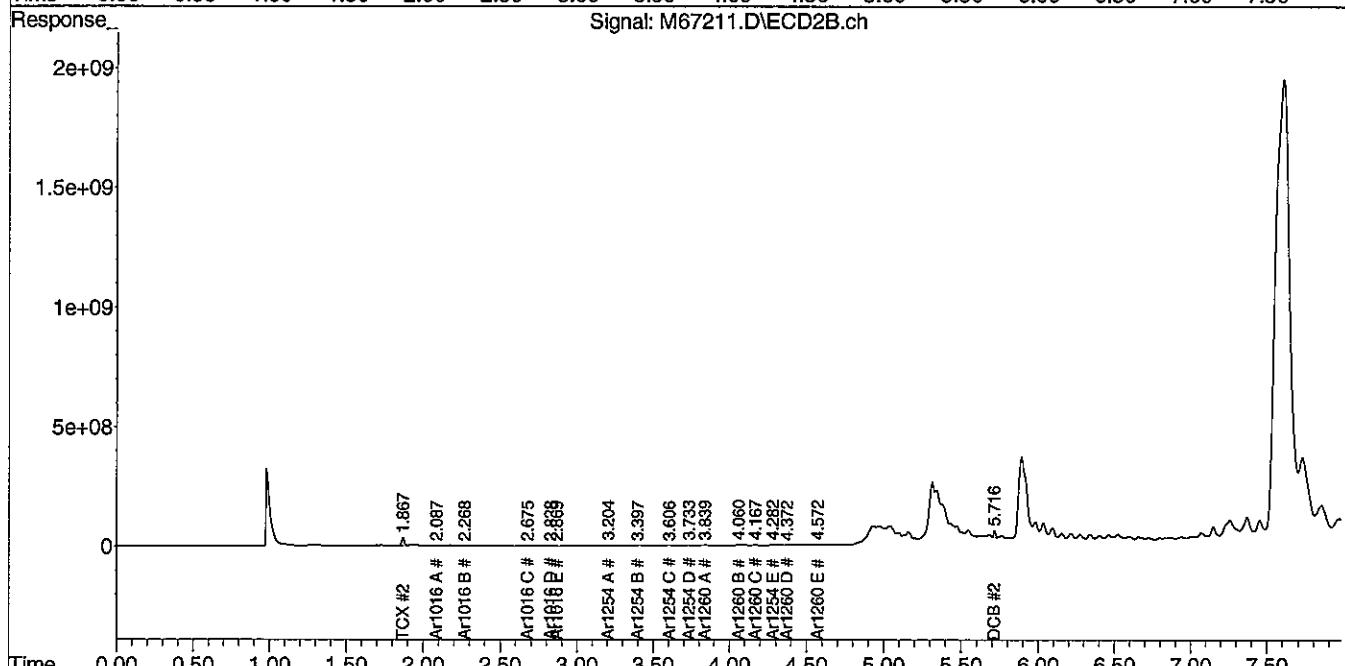
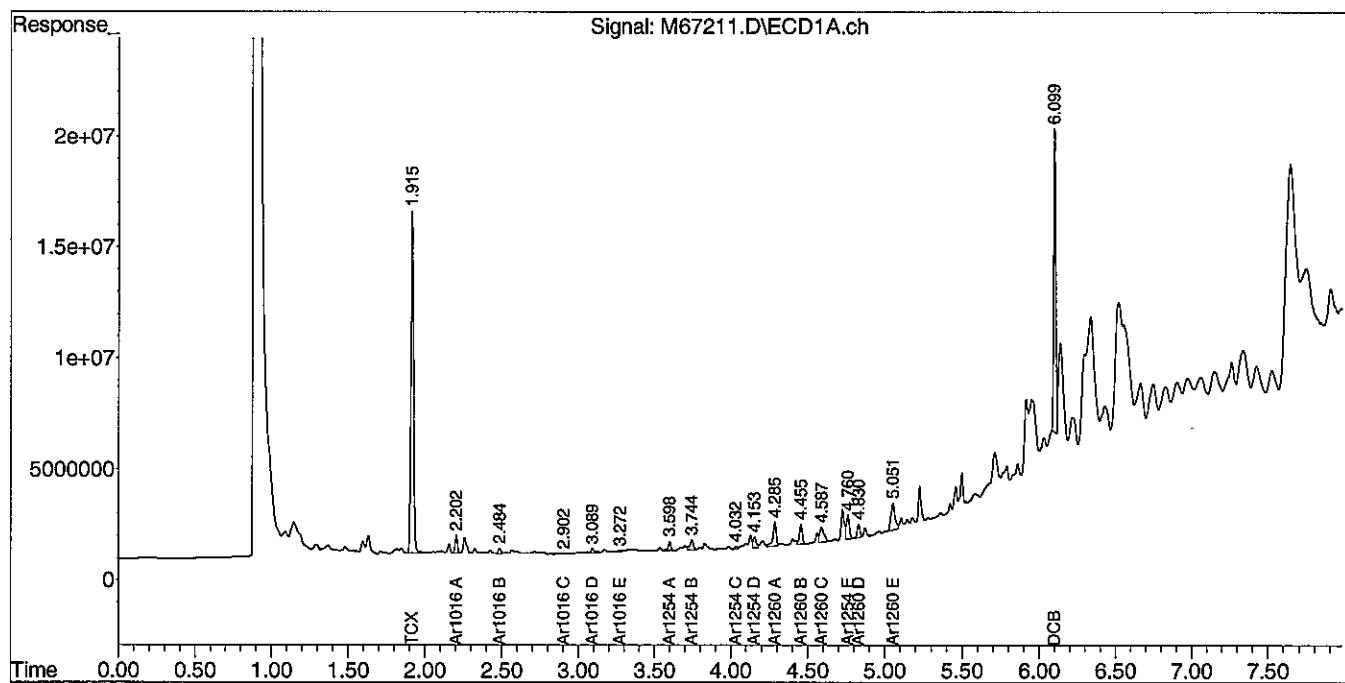
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67211.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 10:46 am
 Operator : JK
 Sample : 74873-1,RR,,A/C
 Misc : SOIL
 ALS Vial : 10 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 16:02:08 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-216 (2-4')

Lab Sample ID: 74873-2
 Matrix: Solid
 Percent Solid: 92
 Dilution Factor: 21
 Collection Date: 02/19/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/20/13
 Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	693	U
PCB-1221	693	U
PCB-1232	693	U
PCB-1242	693	U
PCB-1248	693	U
PCB-1254	693	U
PCB-1260	693	12900

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-2,1:20,,A/C

Column ID: 0.25 mm

Data File: M67214.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 21.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	12942	10661	19.3	

Column to be used to flag RPD values greater than QC limit of 40%

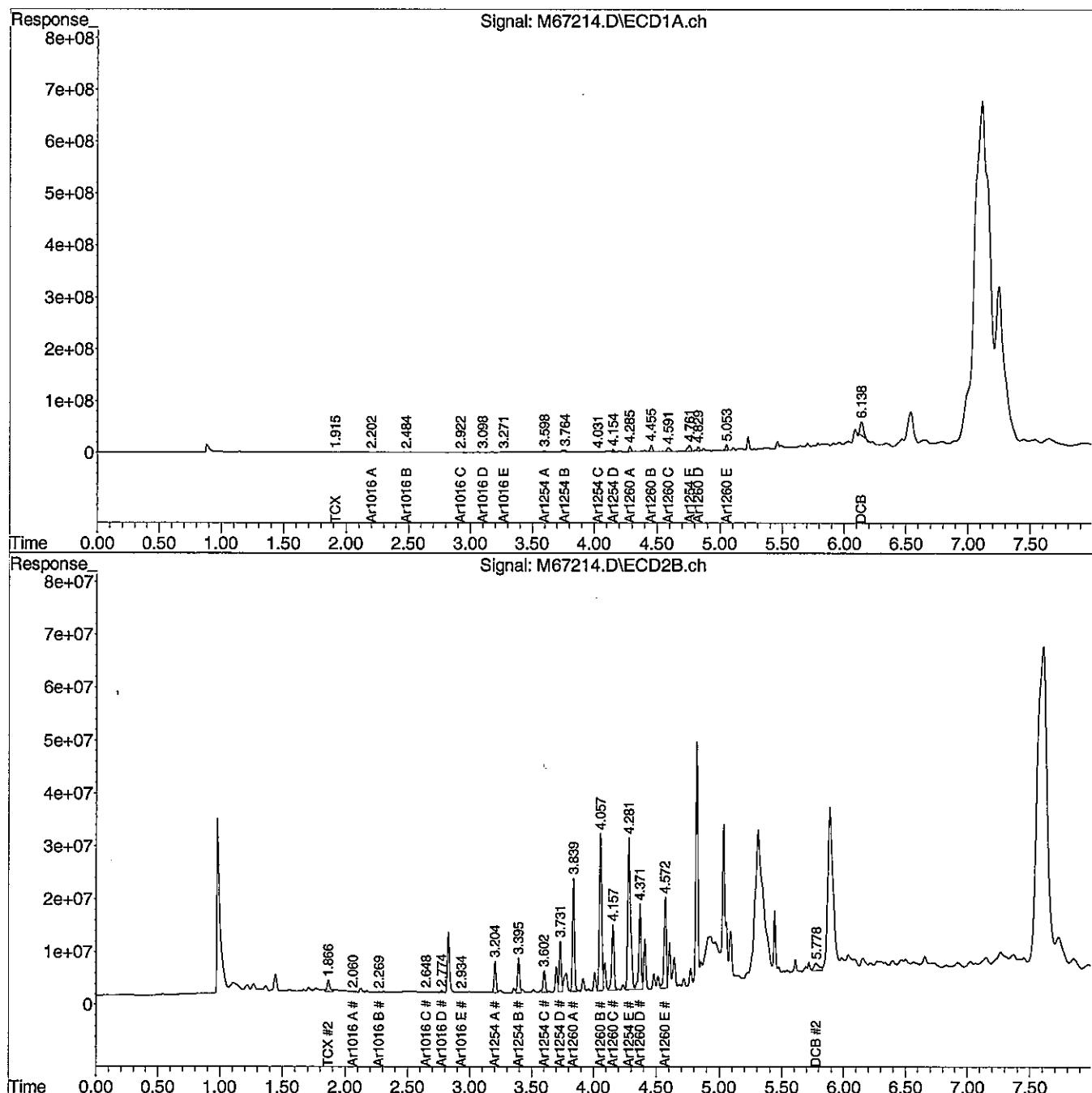
* Values outside QC limits

Comments: _____

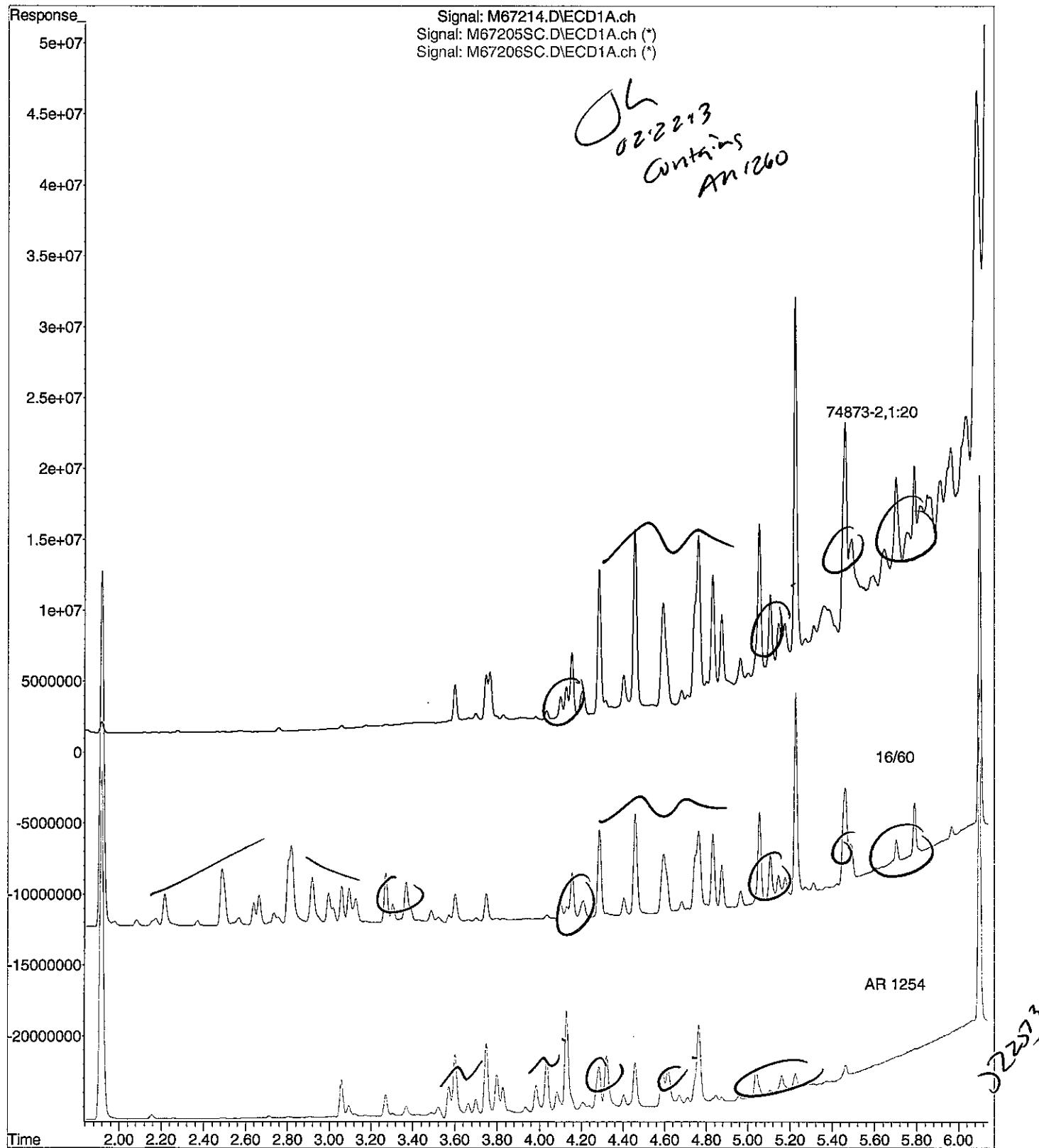
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67214.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 11:16 am
 Operator : JK
 Sample : 74873-2,1:20,,A/C
 Misc : SOIL
 ALS Vial : 13 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:57:46 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67214.D
Operator : JK
Acquired : 22 Feb 2013 11:16 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-2,1:20,,A/C
Misc Info : SOIL
Vial Number: 13



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-216 (4-6')

Lab Sample ID: 74873-3 RR
 Matrix: Solid
 Percent Solid: 92
 Dilution Factor: 1.0
 Collection Date: 02/19/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/20/13
 Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	112	%
Decachlorobiphenyl	76	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

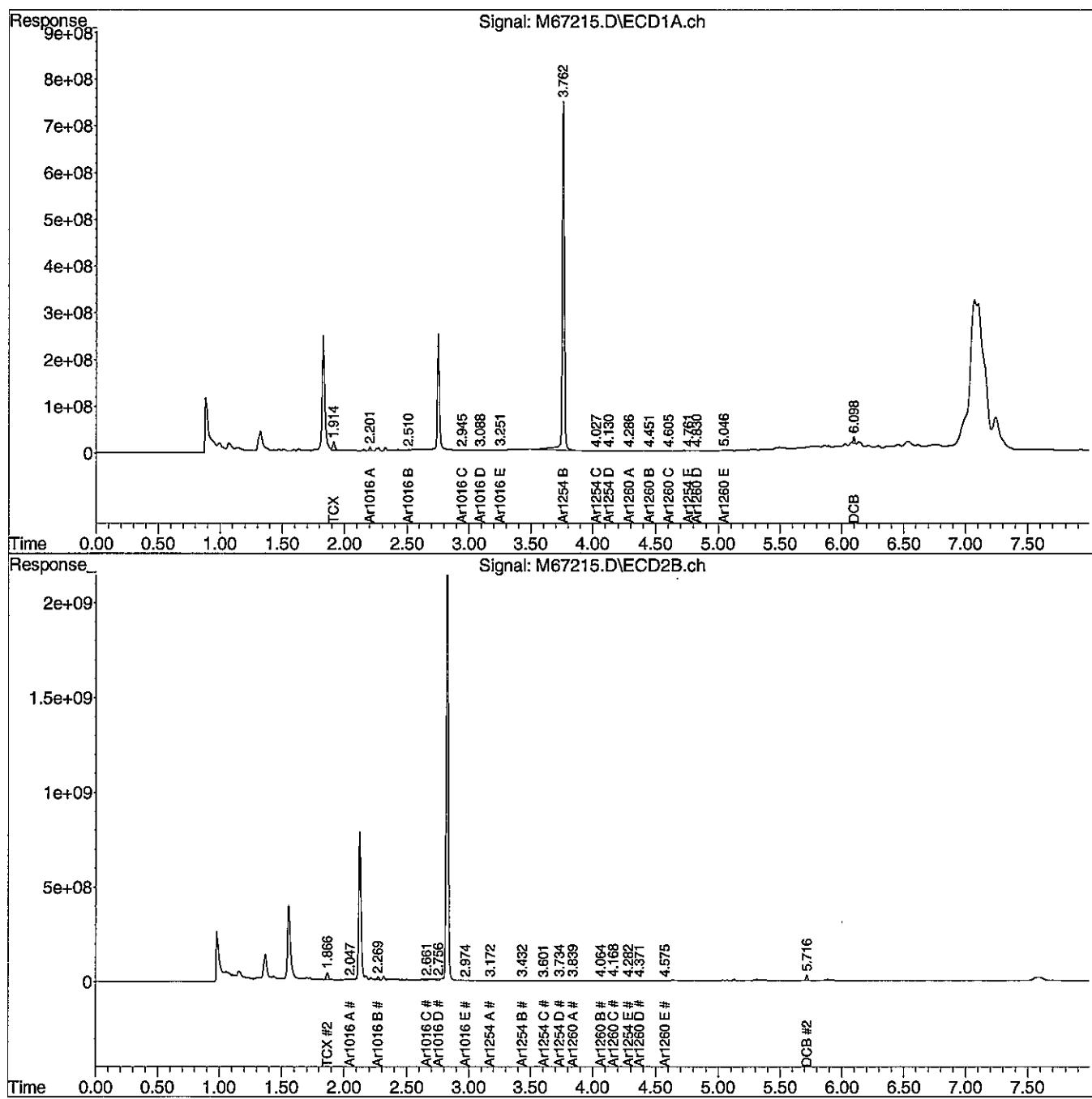
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67215.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 11:27 am
 Operator : JK
 Sample : 74873-3,RR,,A/C
 Misc : SOIL
 ALS Vial : 14 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 16:23:45 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-215 (0-2')

Lab Sample ID: 74873-4
Matrix: Solid
Percent Solid: 97
Dilution Factor: 5.0
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/20/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	102	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

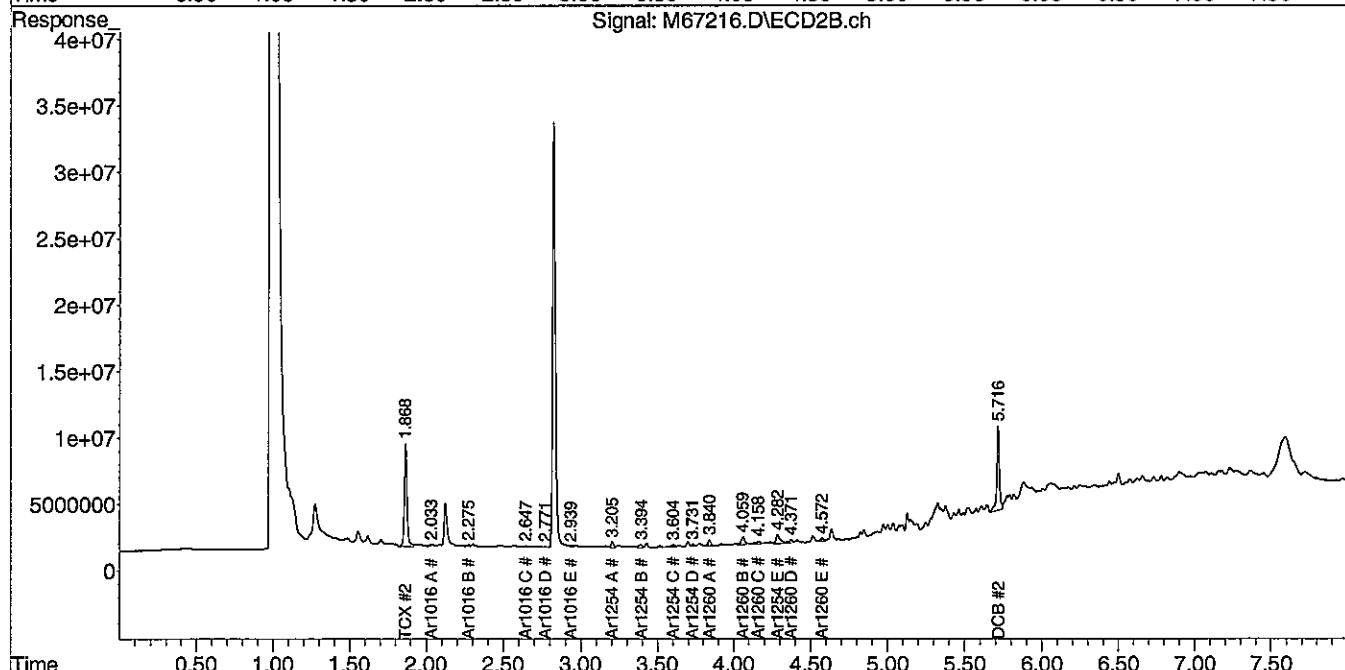
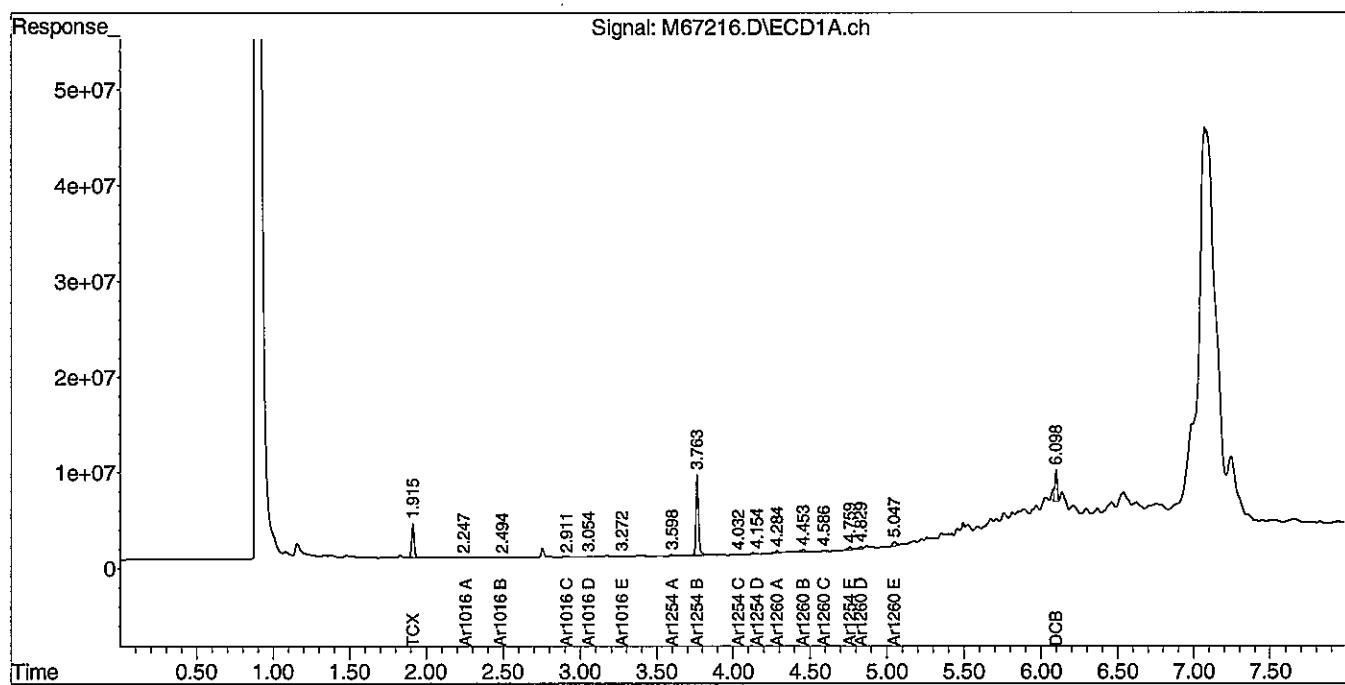
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * Quantitation limits increased due to the sample matrix affect.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67216.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 11:37 am
 Operator : JK
 Sample : 74873-4,1:5,,A/C
 Misc : SOIL
 ALS Vial : 15 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 16:26:09 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-215 (2-4")

Lab Sample ID: 74873-5
Matrix: Solid
Percent Solid: 93
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/20/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	3210

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	84	%
Decachlorobiphenyl	52	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-5,1:5,,A/C

Column ID: 0.25 mm

Data File: M67217.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3213	2856	11.7	

Column to be used to flag RPD values greater than QC limit of 40%

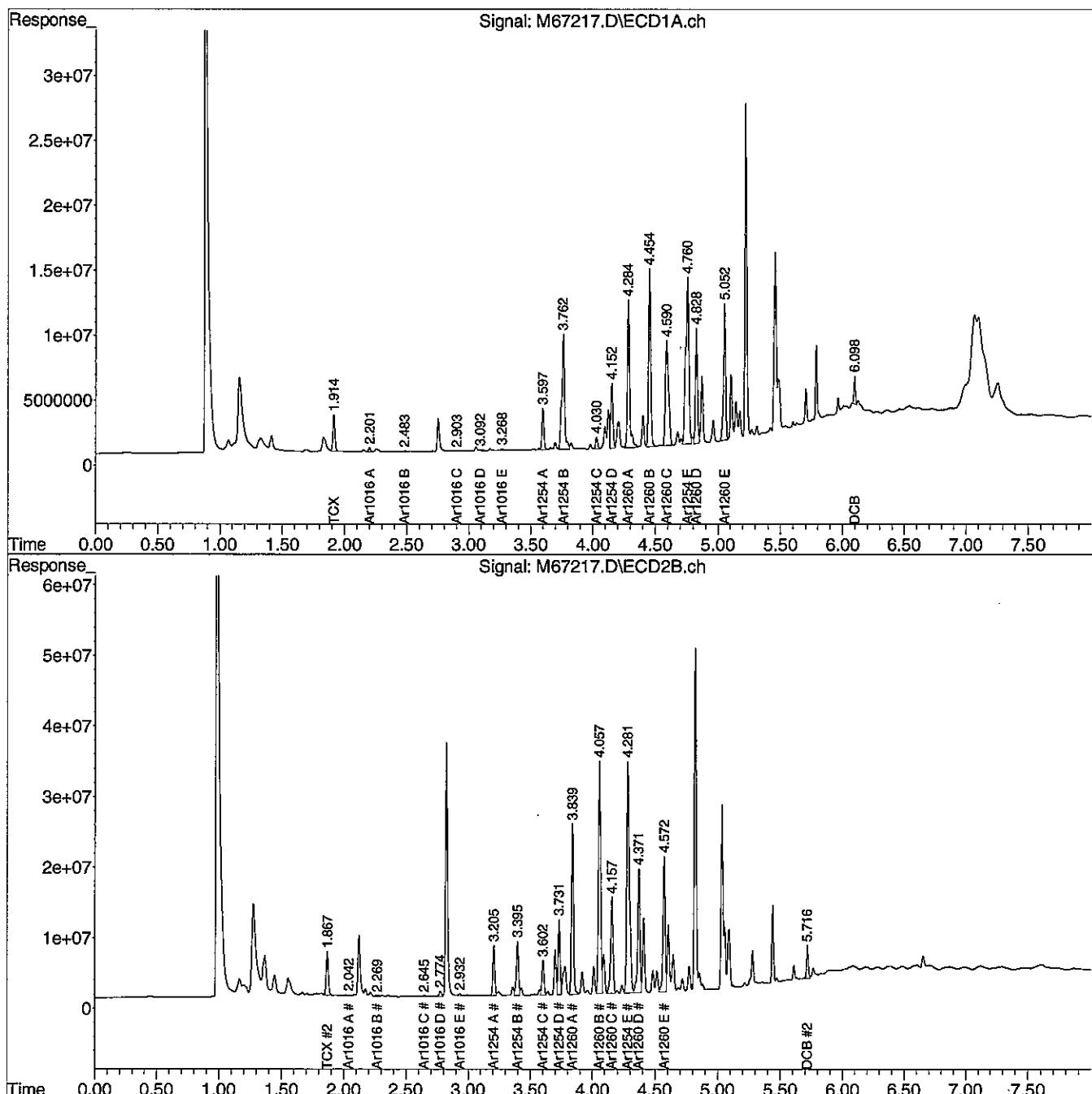
* Values outside QC limits

Comments: _____

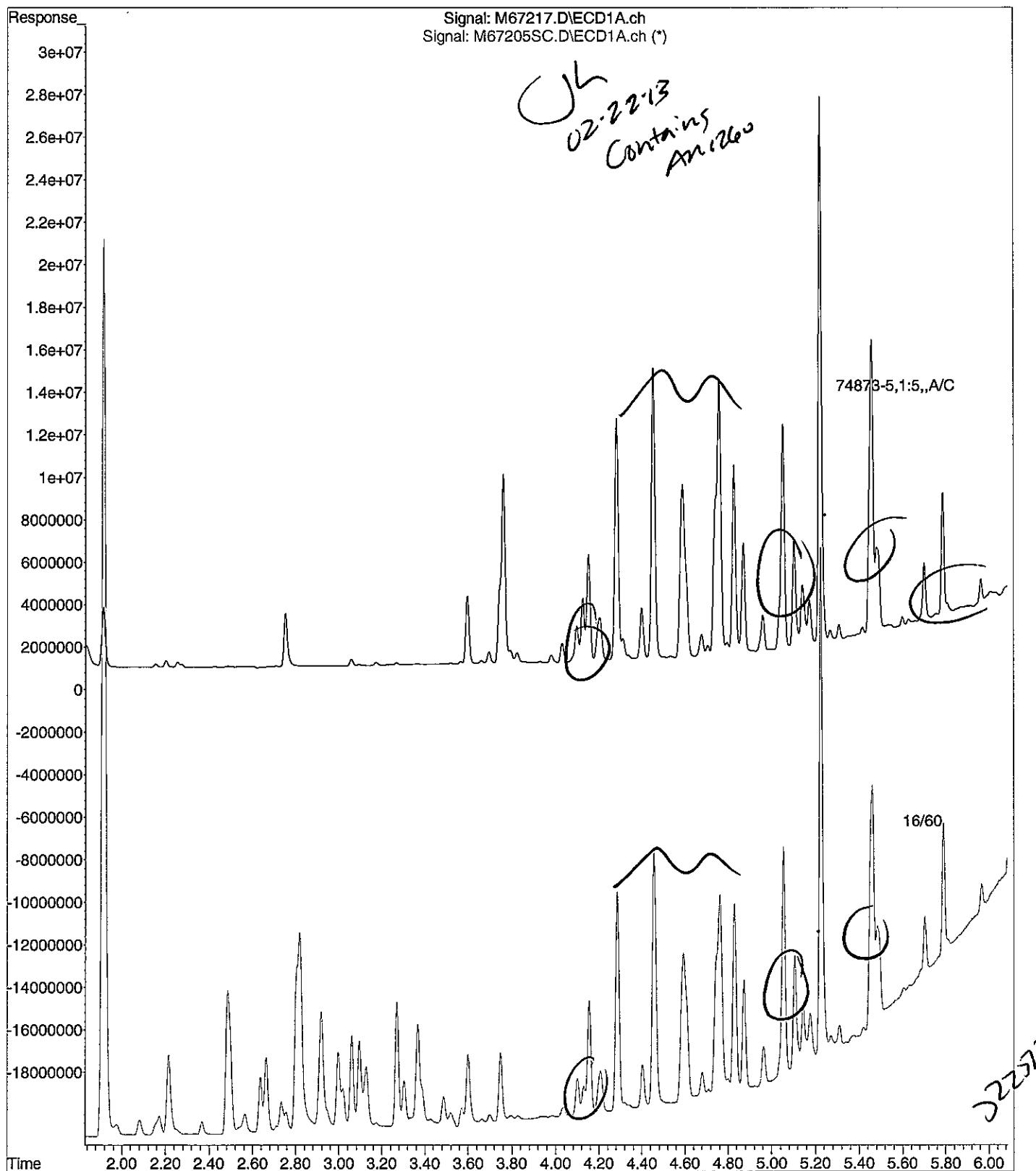
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67217.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 11:47 am
 Operator : JK
 Sample : 74873-5,1:5,,A/C
 Misc : SOIL
 ALS Vial : 16 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:57:52 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67217.D
Operator : JK
Acquired : 22 Feb 2013 11:47 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-5,1:5,,A/C
Misc Info : SOIL
Vial Number: 16



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-215 (4-6')

Lab Sample ID: 74873-6
Matrix: Solid
Percent Solid: 93
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/20/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	2260

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	97	%
Decachlorobiphenyl	61	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-6,1:5,,A/C

Column ID: 0.25 mm

Data File: M67218.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.3

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2264	2006	12.1	

Column to be used to flag RPD values greater than QC limit of 40%

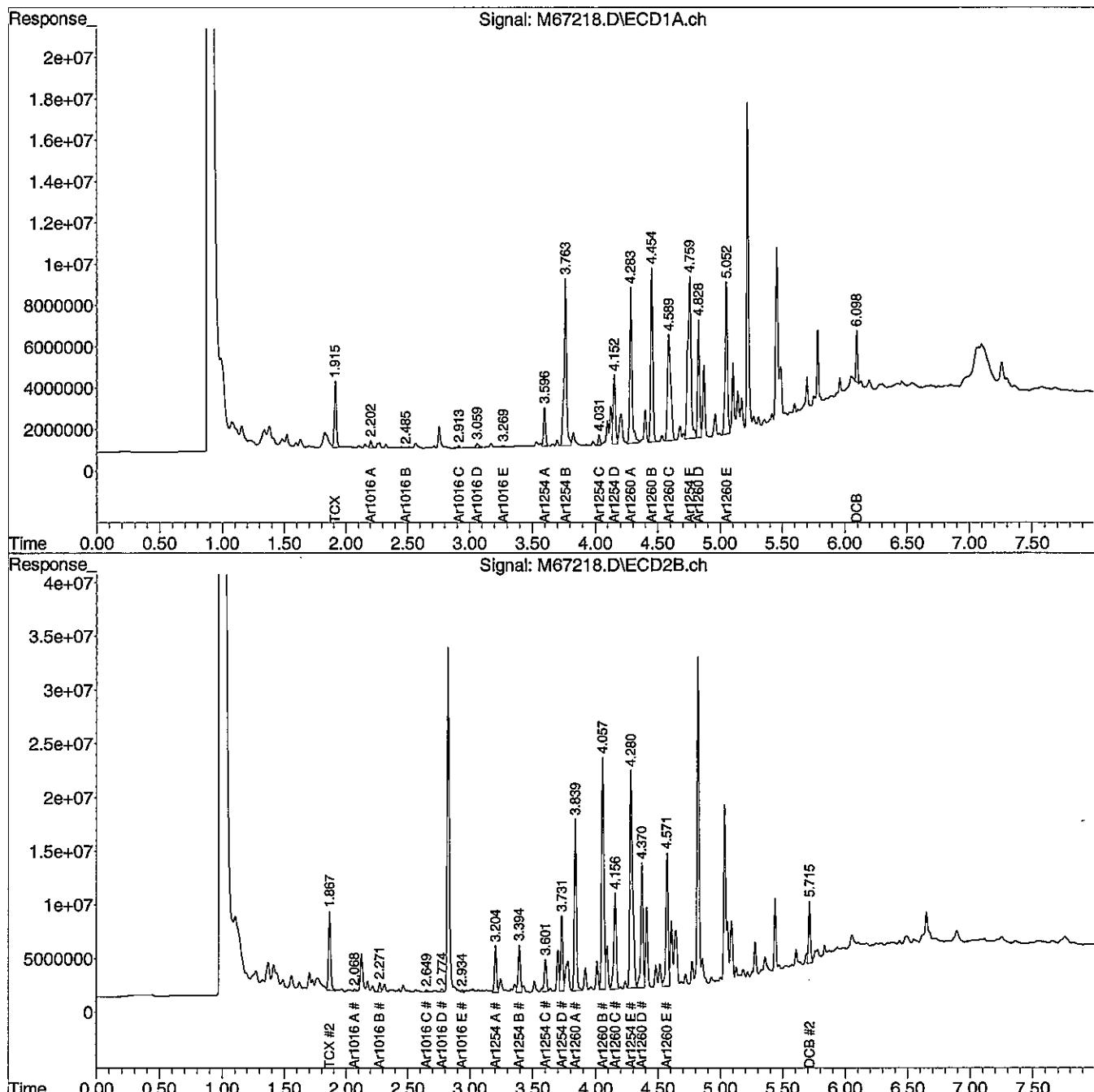
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67218.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 11:57 am
 Operator : JK
 Sample : 74873-6,1:5,,A/C
 Misc : SOIL
 ALS Vial : 17 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:57:54 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-215 (6-8")

Lab Sample ID: 74873-7
Matrix: Solid
Percent Solid: 90
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/20/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	3220

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	87	%
Decachlorobiphenyl	59	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-7,1:5,,A/C

Column ID: 0.25 mm

Data File: M67219.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3220	2722	16.7	

Column to be used to flag RPD values greater than QC limit of 40%

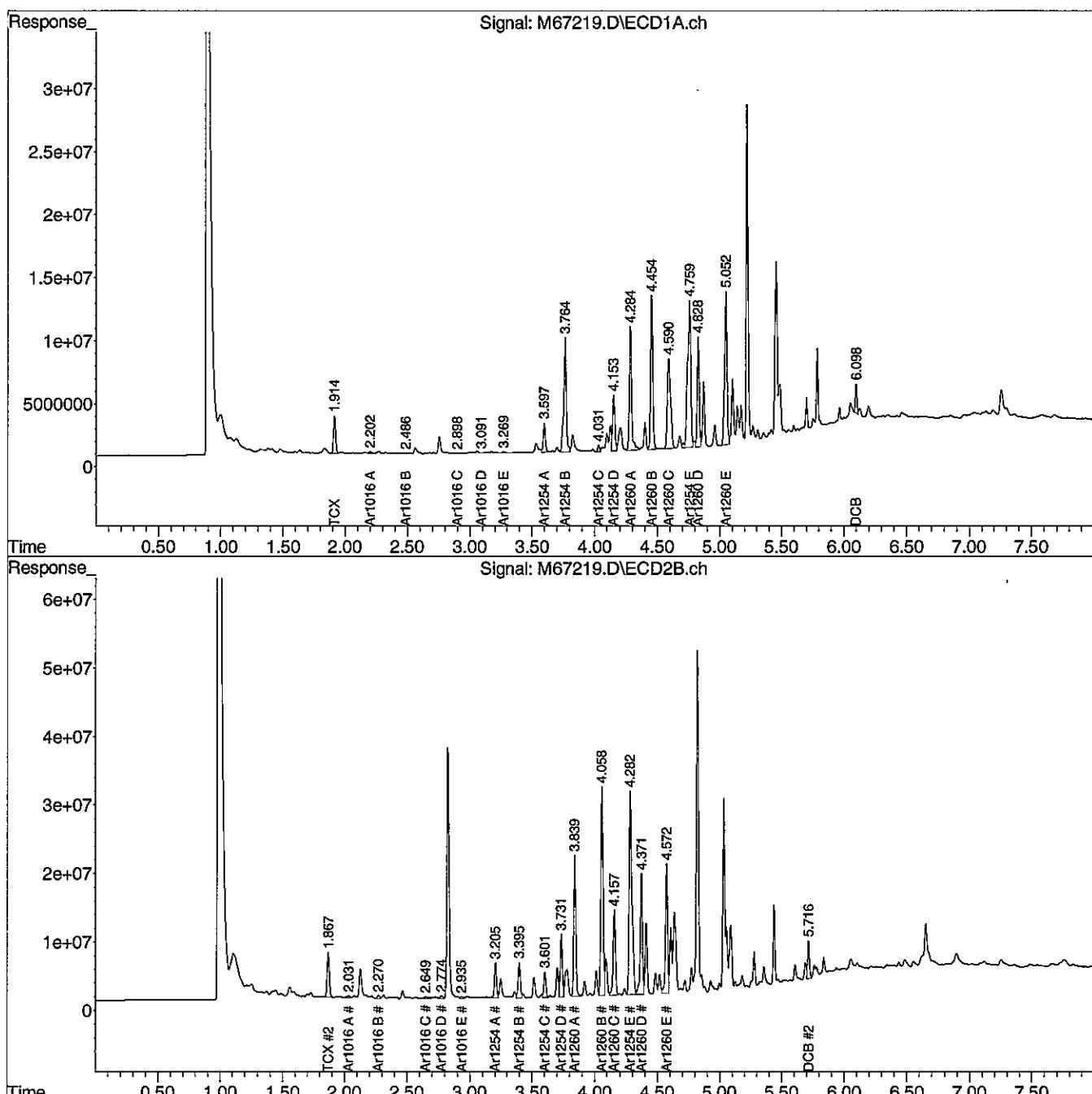
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67219.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 12:07 pm
 Operator : JK
 Sample : 74873-7,1:5,,A/C
 Misc : SOIL
 ALS Vial : 18 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:57:56 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 22, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-215 (8-10')

Lab Sample ID: 74873-8
Matrix: Solid
Percent Solid: 88
Dilution Factor: 54
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/20/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	1780	U
PCB-1221	1780	U
PCB-1232	1780	U
PCB-1242	1780	U
PCB-1248	1780	U
PCB-1254	1780	U
PCB-1260	1780	44600

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-8,1:50,,A/C

Column ID: 0.25 mm

Data File: M67220.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 54.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	44645	40316	10.2	

Column to be used to flag RPD values greater than QC limit of 40%

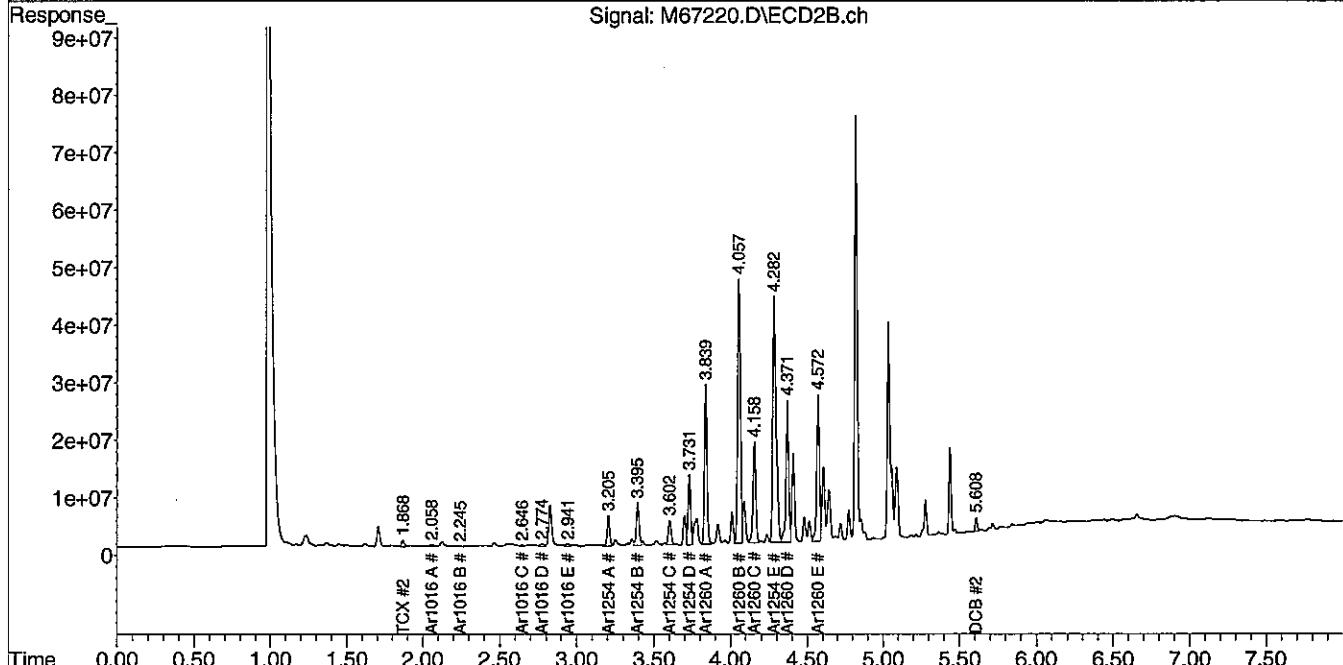
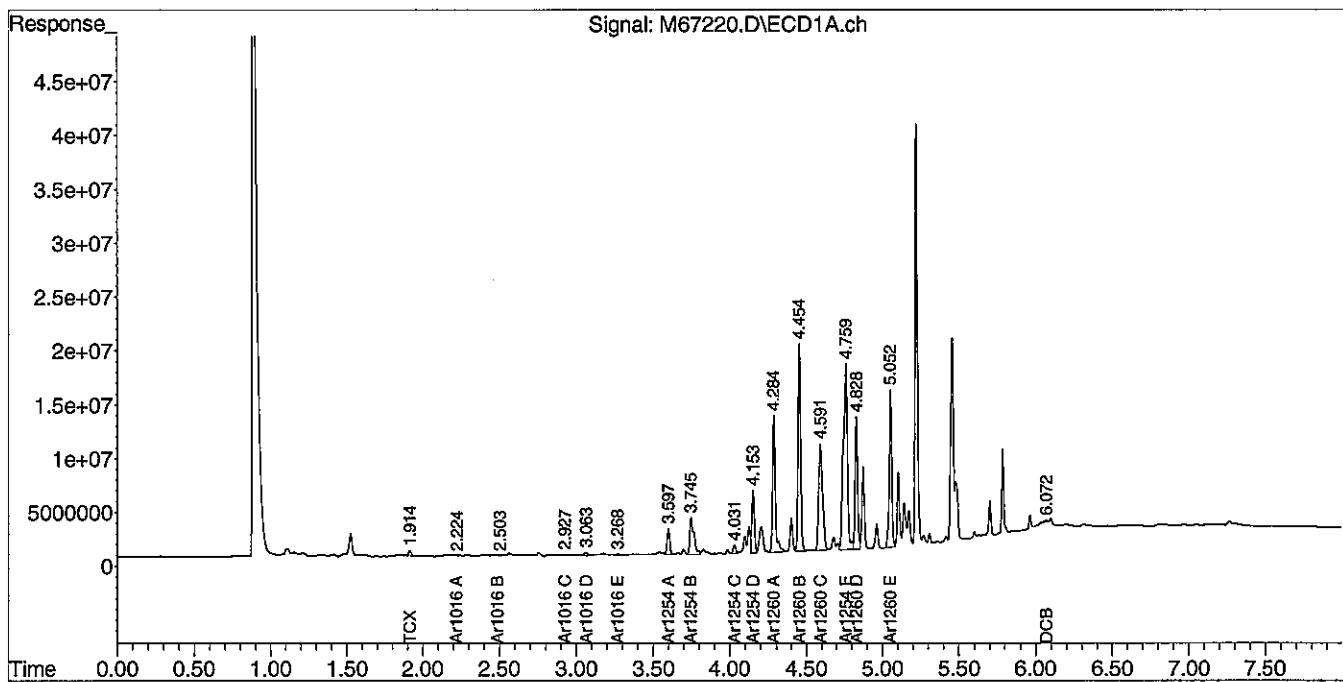
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67220.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 12:17 pm
 Operator : JK
 Sample : 74873-8,1:50,,A/C
 Misc : SOIL
 ALS Vial : 19 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:57:58 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-214 (0-2')

Lab Sample ID: 74873-9 RX
Matrix: Solid
Percent Solid: 88
Dilution Factor: 6
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/25/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	385

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	115	%
Decachlorobiphenyl	114	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-9,RX,1:5

Column ID: 0.25 mm

Data File: M67419.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	385	343	11.6	

Column to be used to flag RPD values greater than QC limit of 40%

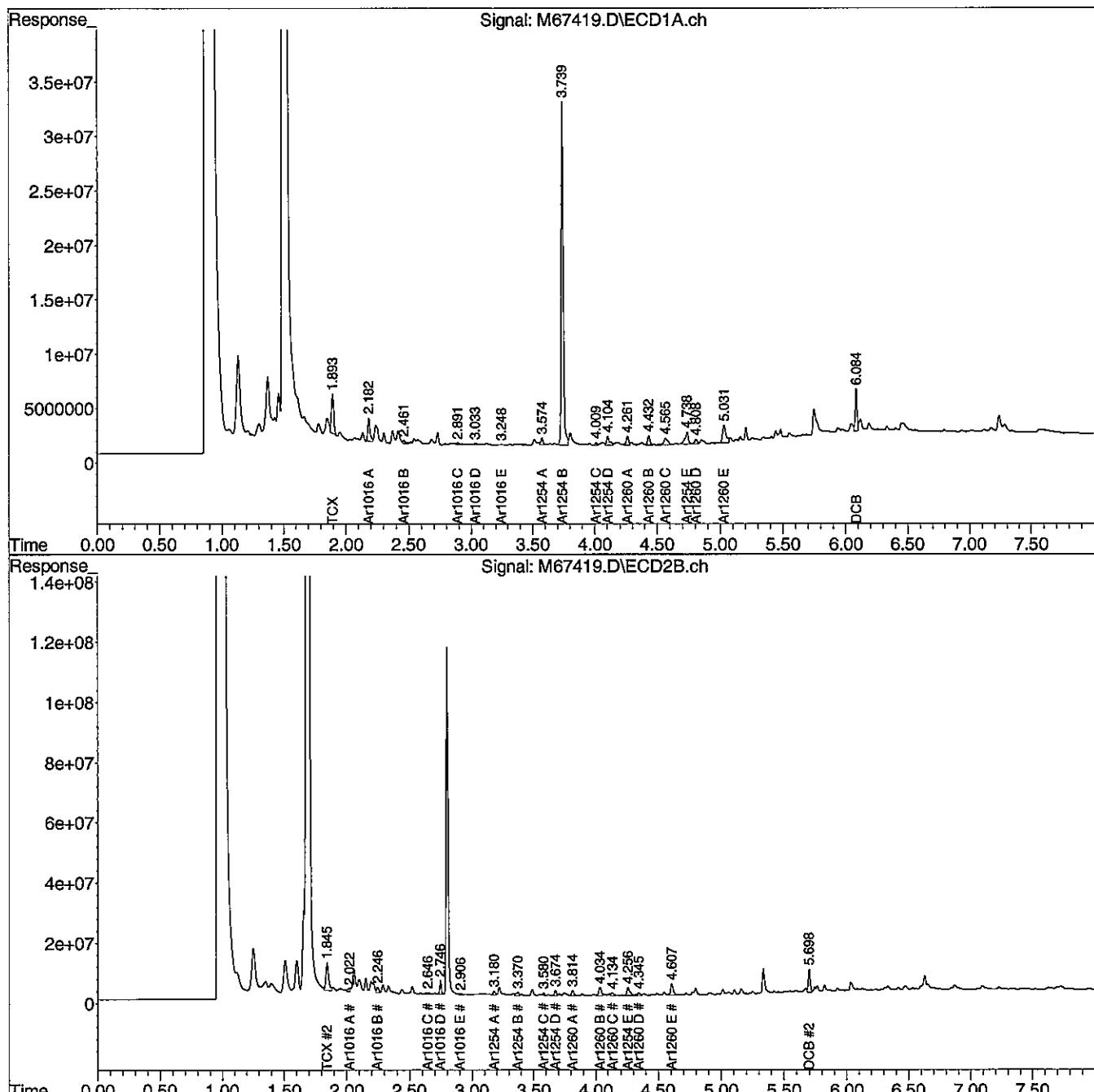
* Values outside QC limits

Comments: _____

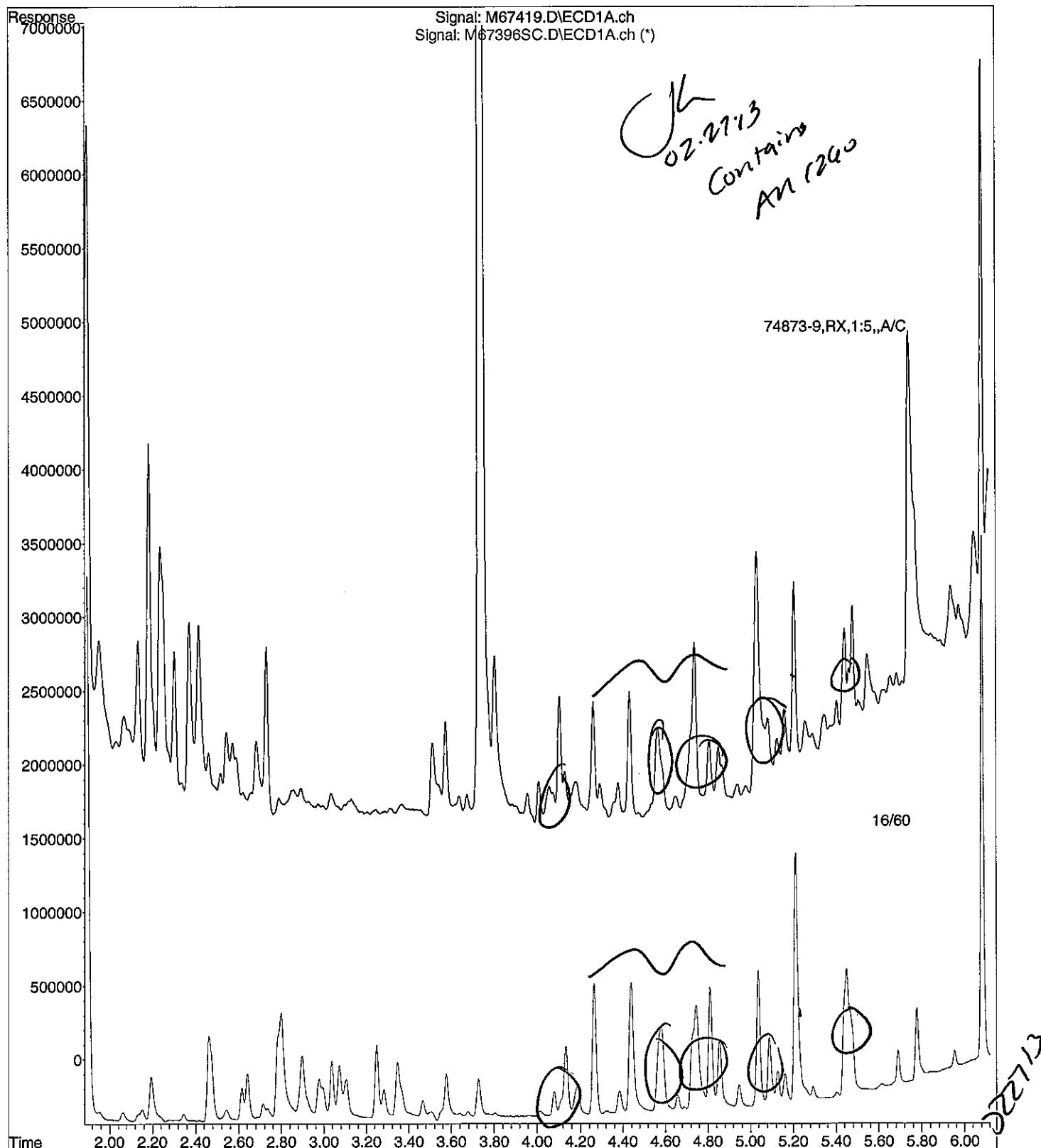
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67419.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 7:53 pm
 Operator : JK
 Sample : 74873-9, RX, 1:5
 Misc : SOIL
 ALS Vial : 11 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:44:18 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67419.D
Operator : JK
Acquired : 26 Feb 2013 7:53 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-9,RX,1:5
Misc Info : SOIL
Vial Number: 11



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-214 (2-4')

Lab Sample ID: 74873-10 RX
Matrix: Solid
Percent Solid: 91
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/25/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	1280

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	67	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-10,RX,1:5

Column ID: 0.25 mm

Data File: M67420.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1275	942	30.0	

Column to be used to flag RPD values greater than QC limit of 40%

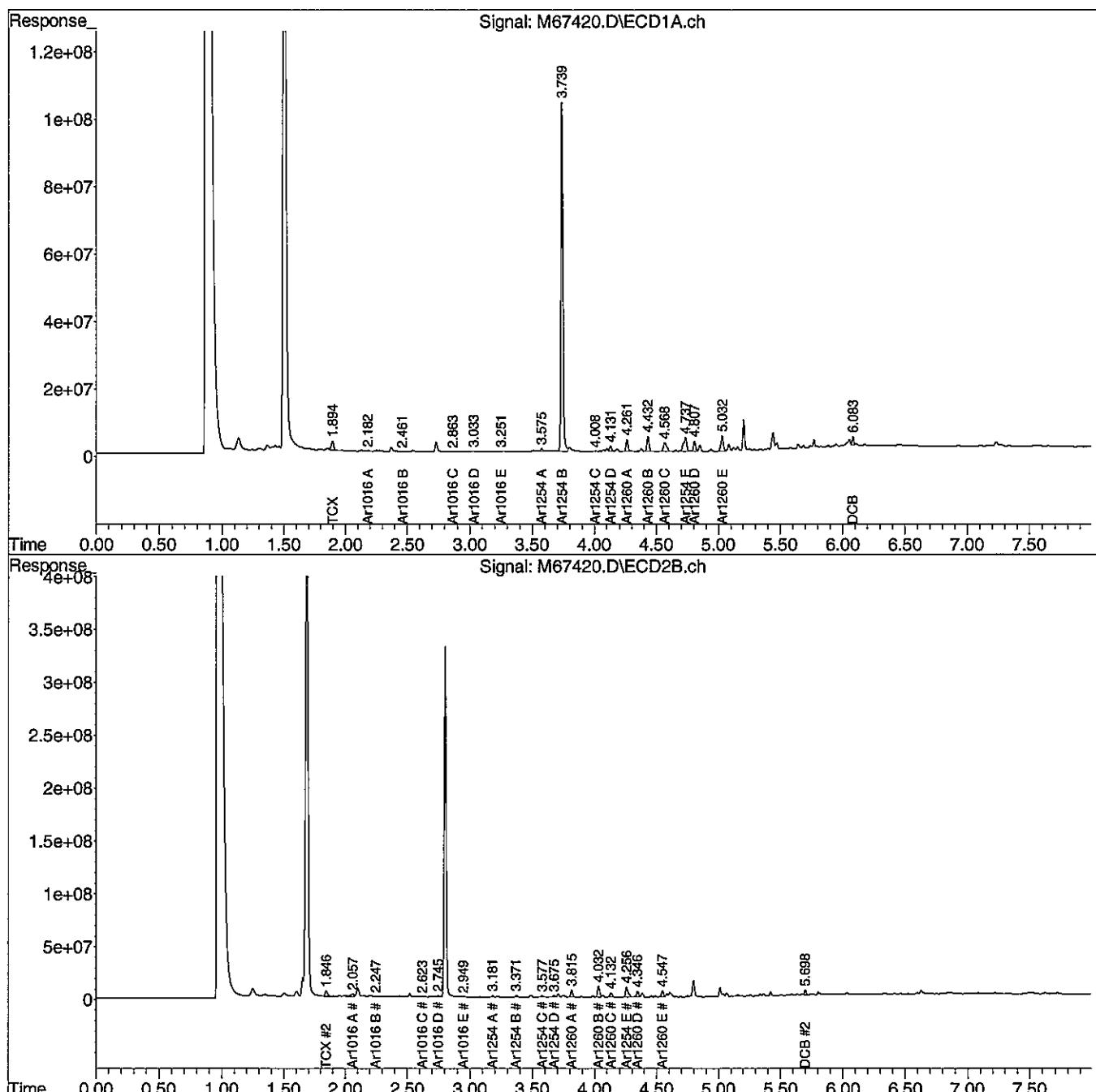
* Values outside QC limits

Comments: _____

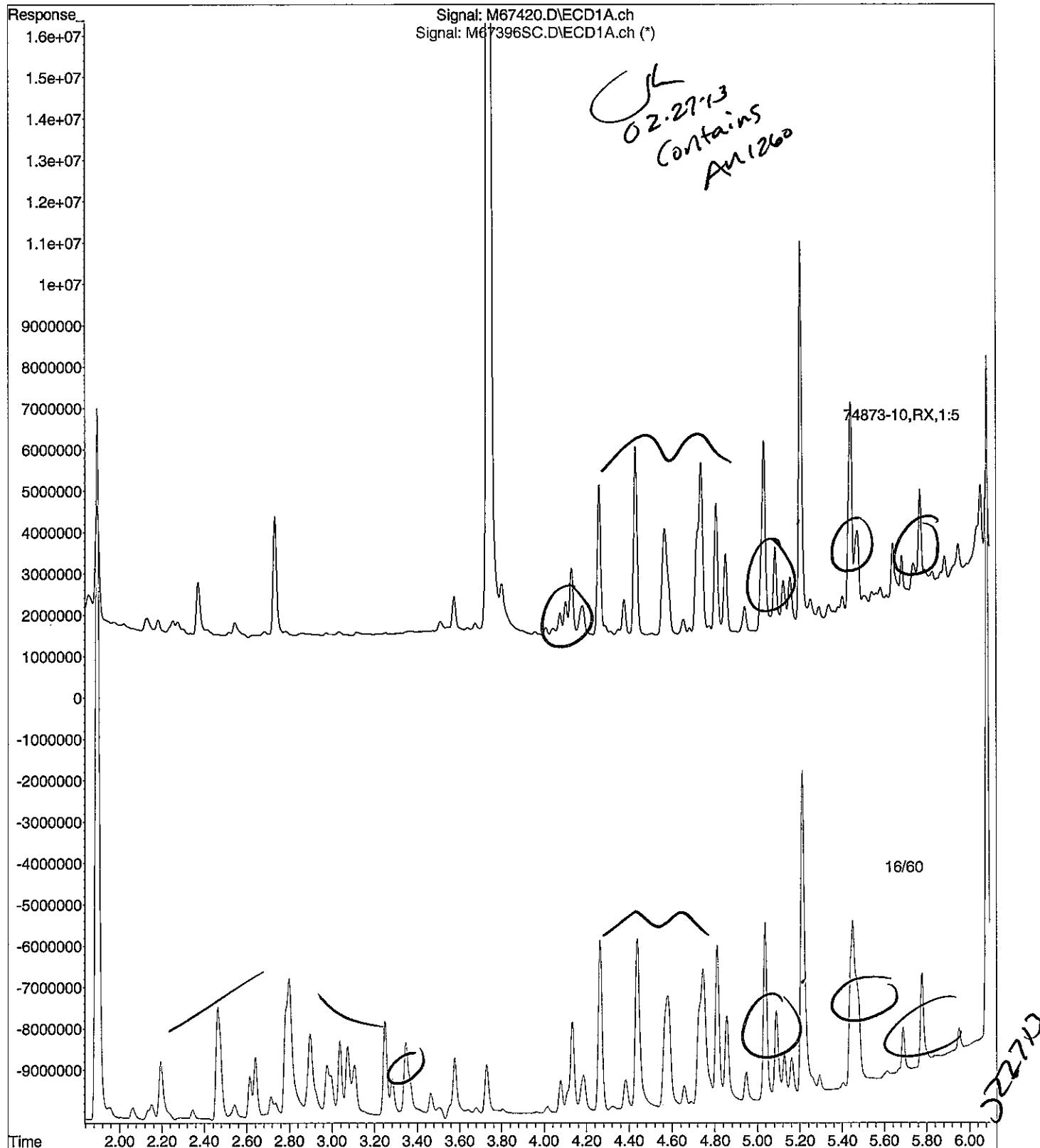
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67420.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 8:03 pm
 Operator : JK
 Sample : 74873-10,RX,1:5
 Misc : SOIL
 ALS Vial : 12 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:45:28 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67420.D
Operator : JK
Acquired : 26 Feb 2013 8:03 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-10,RX,1:5
Misc Info : SOIL
Vial Number: 12



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-214 (4-6')

Lab Sample ID: 74873-11
Matrix: Solid
Percent Solid: 88
Dilution Factor: 23
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	759	U
PCB-1221	759	U
PCB-1232	759	U
PCB-1242	759	U
PCB-1248	759	U
PCB-1254	759	U
PCB-1260	759	17200

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-11,1:20,,A/C

Column ID: 0.25 mm

Data File: M67421.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 22.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	17238	14457	17.5	

Column to be used to flag RPD values greater than QC limit of 40%

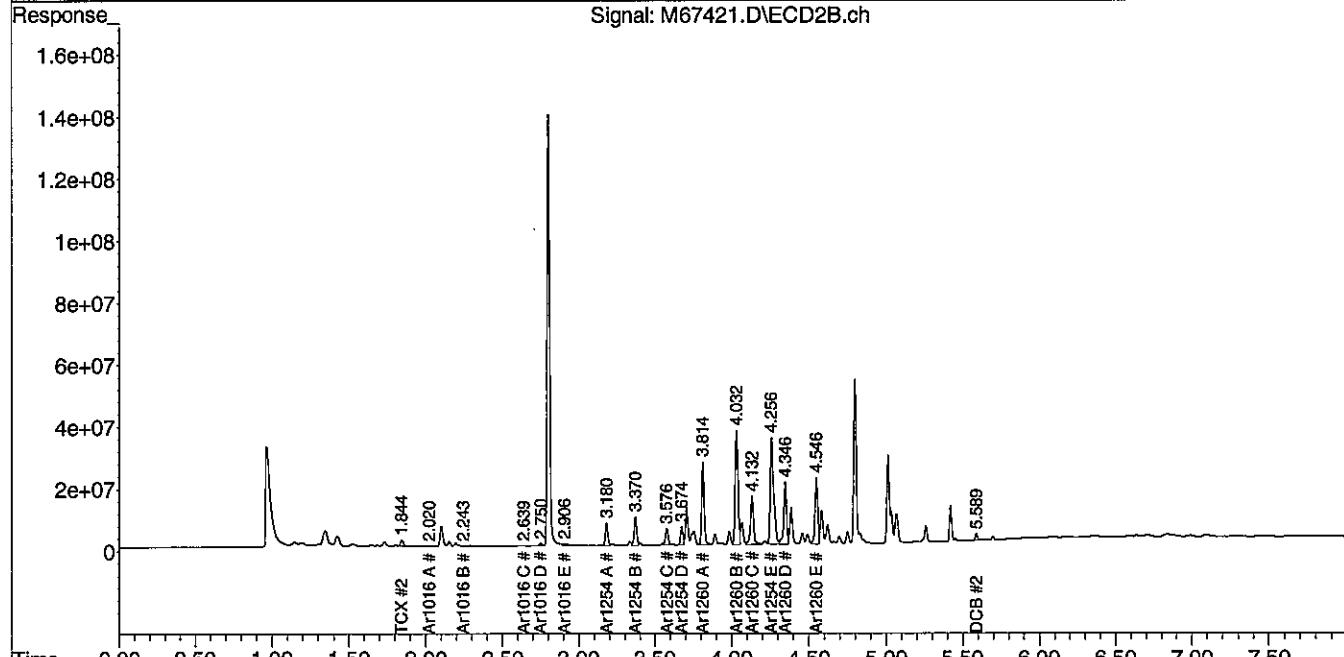
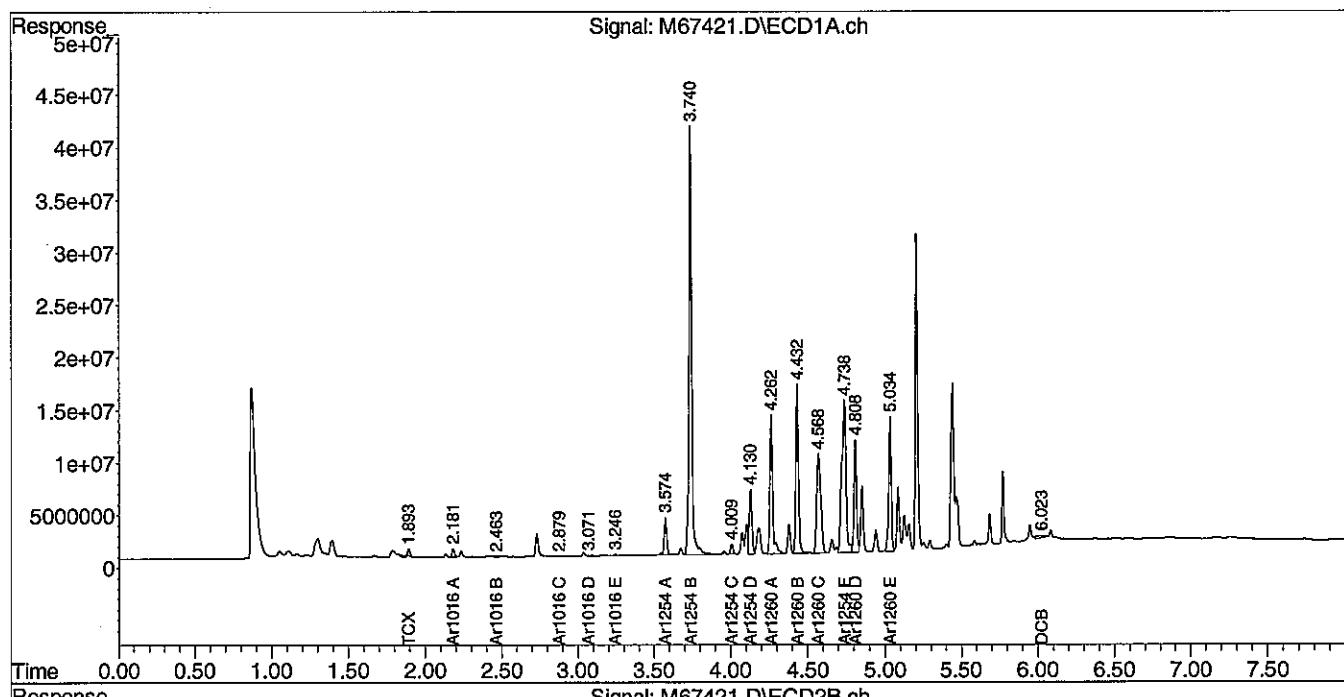
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67421.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 8:14 pm
 Operator : JK
 Sample : 74873-11,1:20,,A/C
 Misc : SOIL
 ALS Vial : 13 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:23:13 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-214 (6-7")

Lab Sample ID: 74873-12
Matrix: Solid
Percent Solid: 87
Dilution Factor: 11
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	8160

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	93	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-12,1:10,,A/C

Column ID: 0.25 mm

Data File: M67422.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.9

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	8161	6808	18.1	

Column to be used to flag RPD values greater than QC limit of 40%

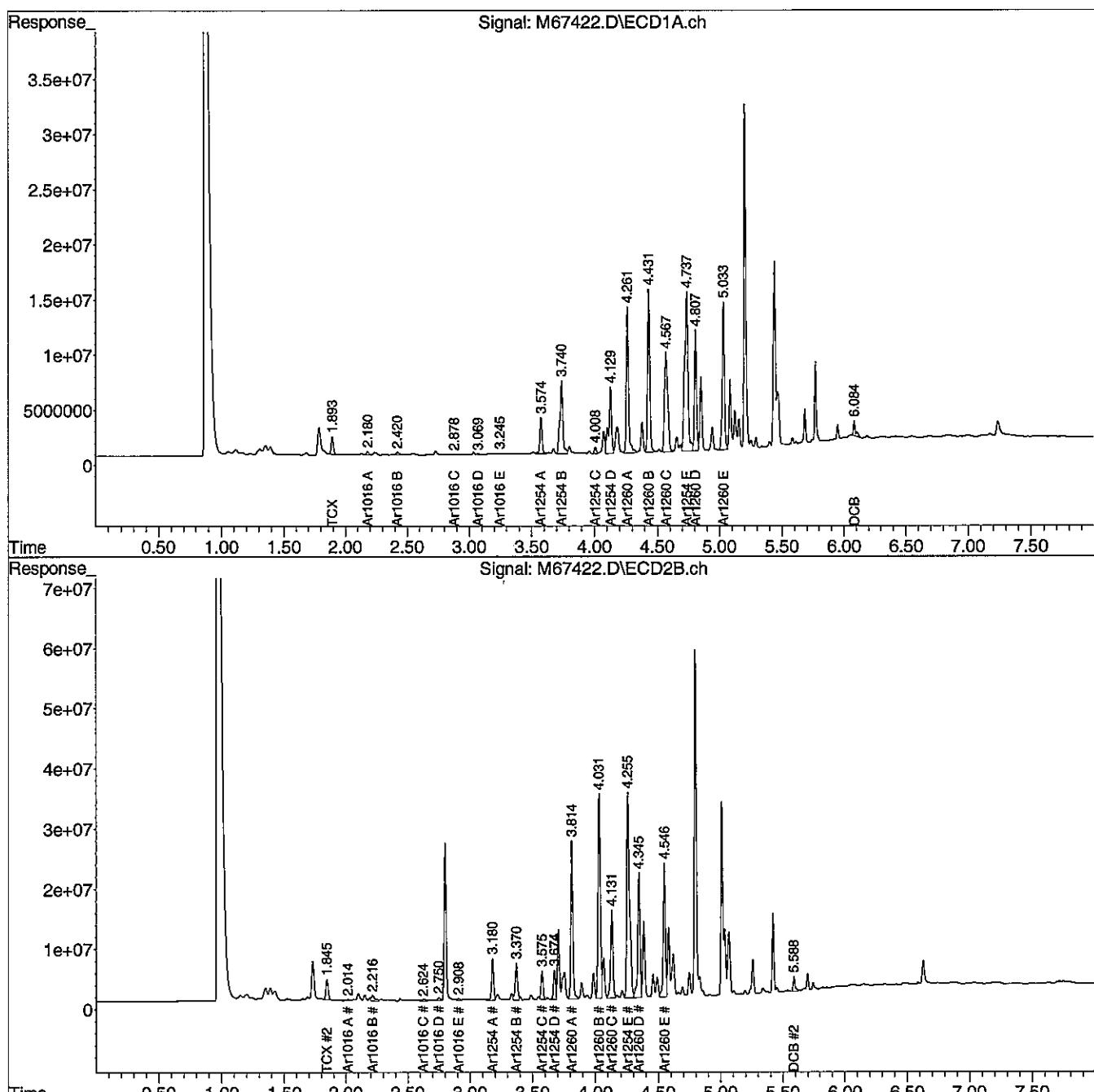
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67422.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 8:24 pm
 Operator : JK
 Sample : 74873-12,1:10,,A/C
 Misc : SOIL
 ALS Vial : 14 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:47:33 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 26, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-213 (0-2')

Lab Sample ID: 74873-13
Matrix: Solid
Percent Solid: 91
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	707

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	111	%
Decachlorobiphenyl	80	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-13,1:5,,A/C

Column ID: 0.25 mm

Data File: M67264.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.3

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	707	625	12.3	

Column to be used to flag RPD values greater than QC limit of 40%

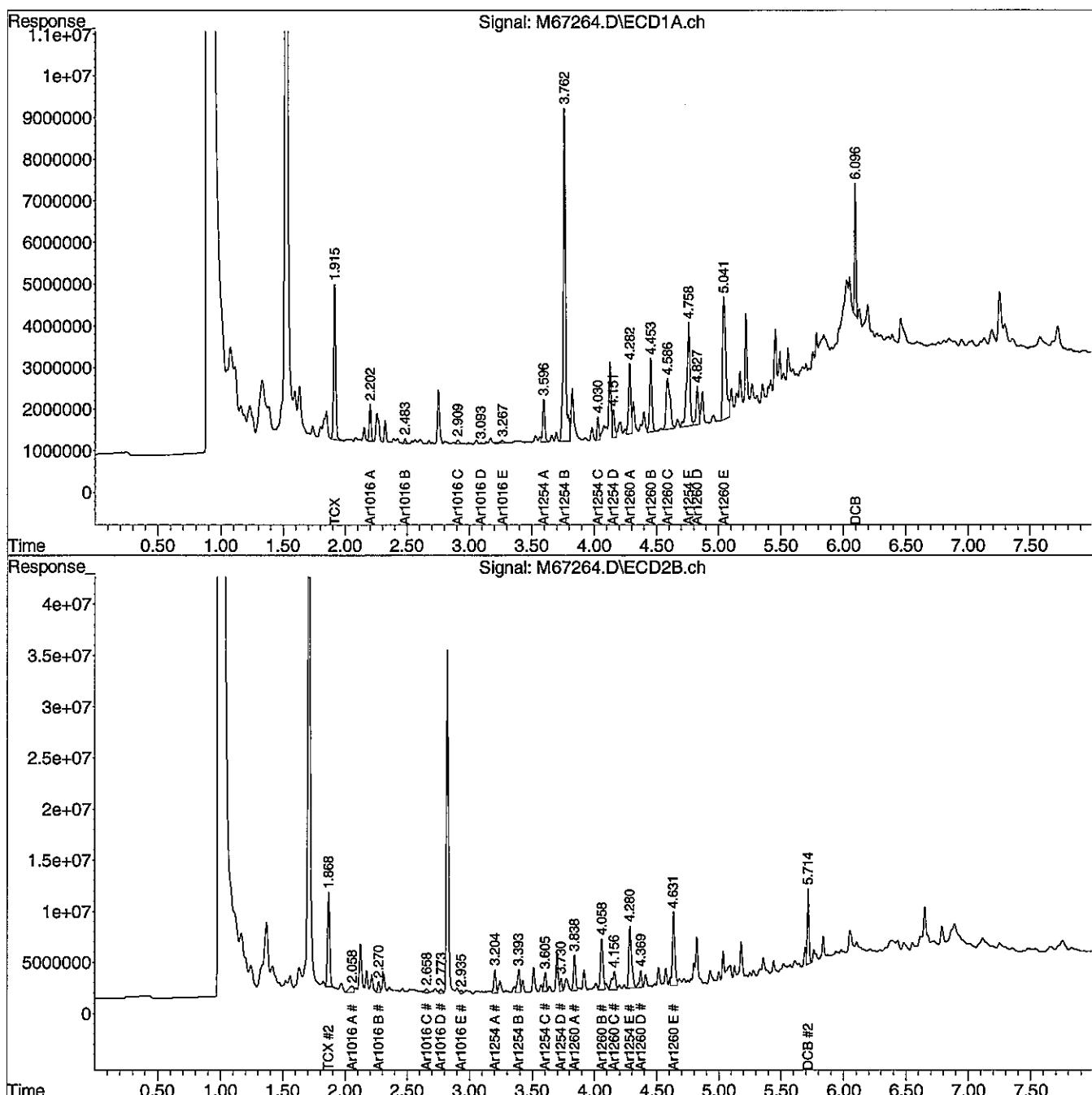
* Values outside QC limits

Comments: _____

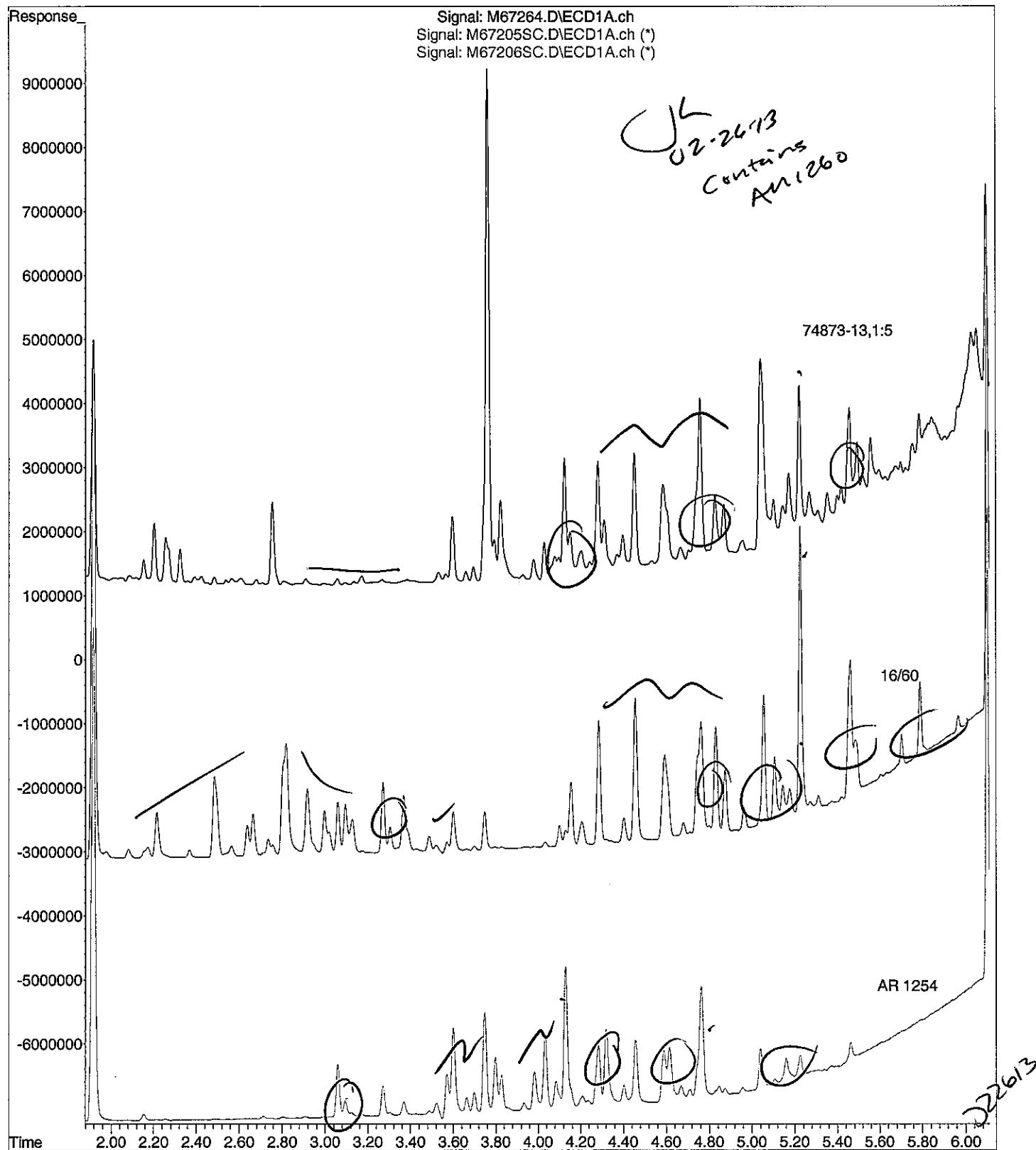
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67264.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 9:00 pm
 Operator : JK
 Sample : 74873-13,1:5,,A/C
 Misc : SOIL
 ALS Vial : 17 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 12:49:20 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67264.D
Operator : JK
Acquired : 22 Feb 2013 9:00 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-13,1:5,,A/C
Misc Info : SOIL
Vial Number: 17



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February 26, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-213 (2-4')

Lab Sample ID: 74873-14
 Matrix: Solid
 Percent Solid: 94
 Dilution Factor: 5
 Collection Date: 02/19/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/21/13
 Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	585

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	78	%
Decachlorobiphenyl	56	%

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-14,1:5,,A/C

Column ID: 0.25 mm

Data File: M67265.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	585	548	6.4	

Column to be used to flag RPD values greater than QC limit of 40%

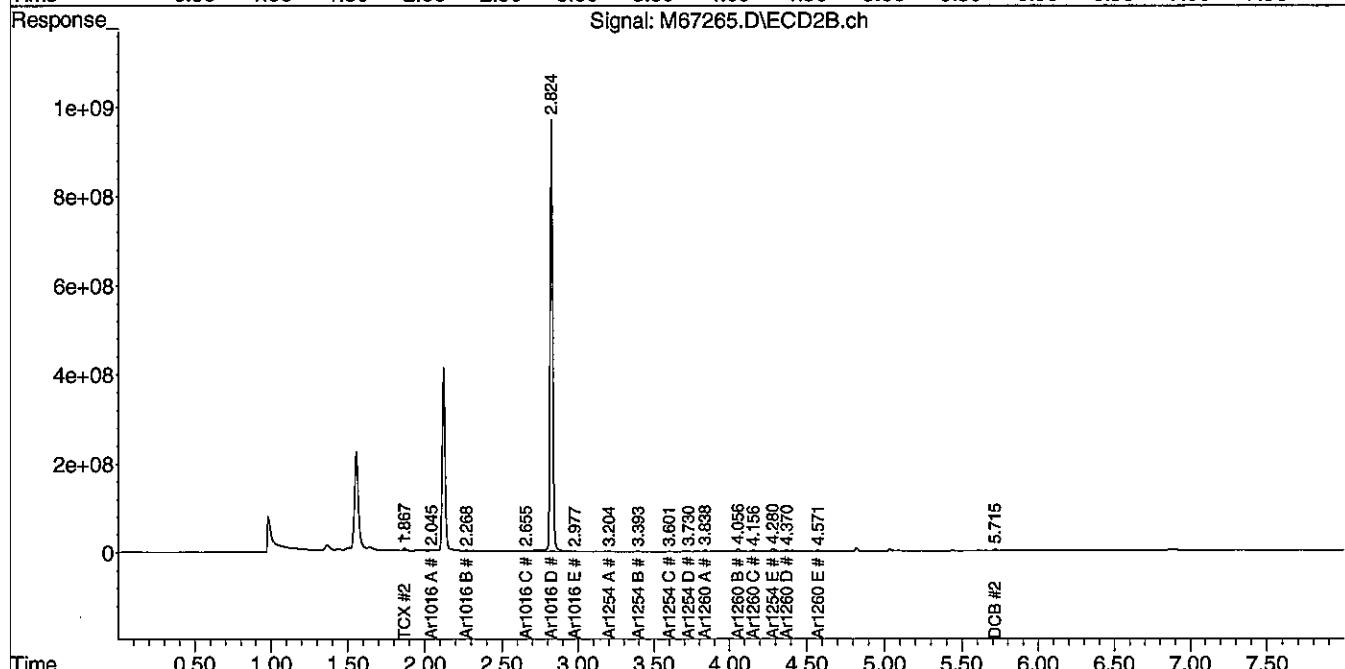
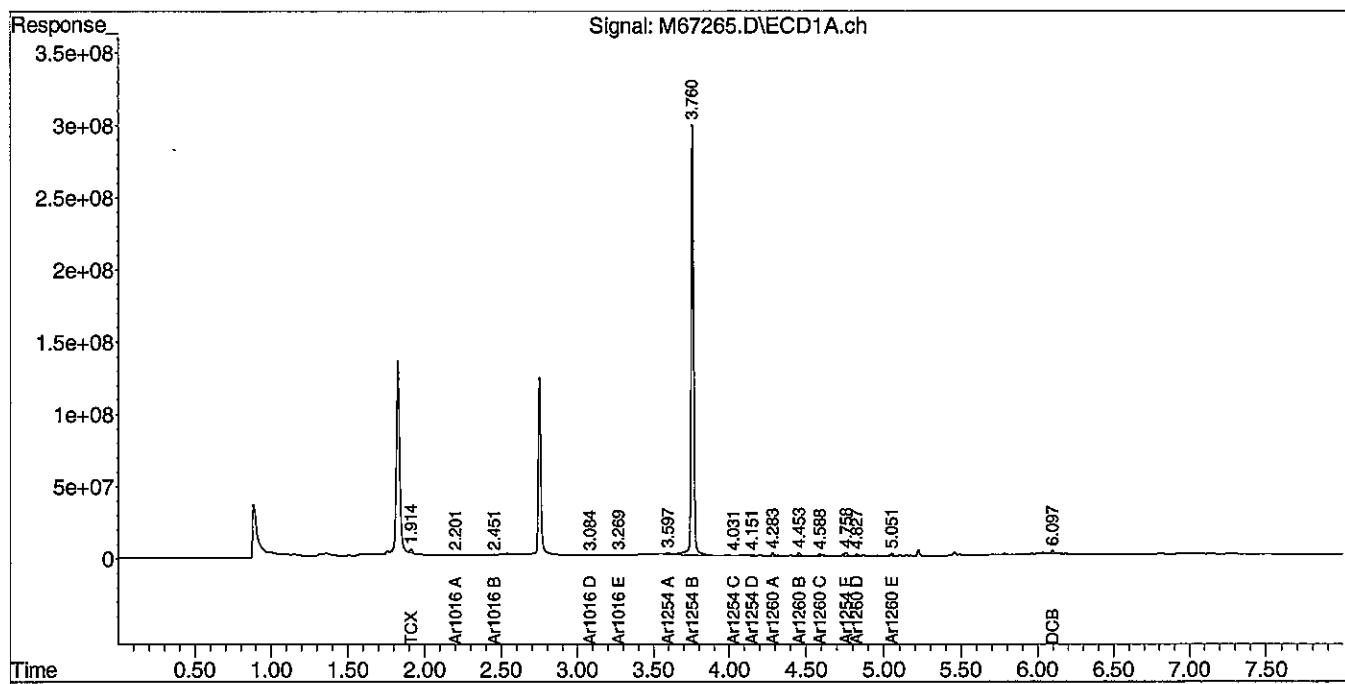
* Values outside QC limits

Comments: _____

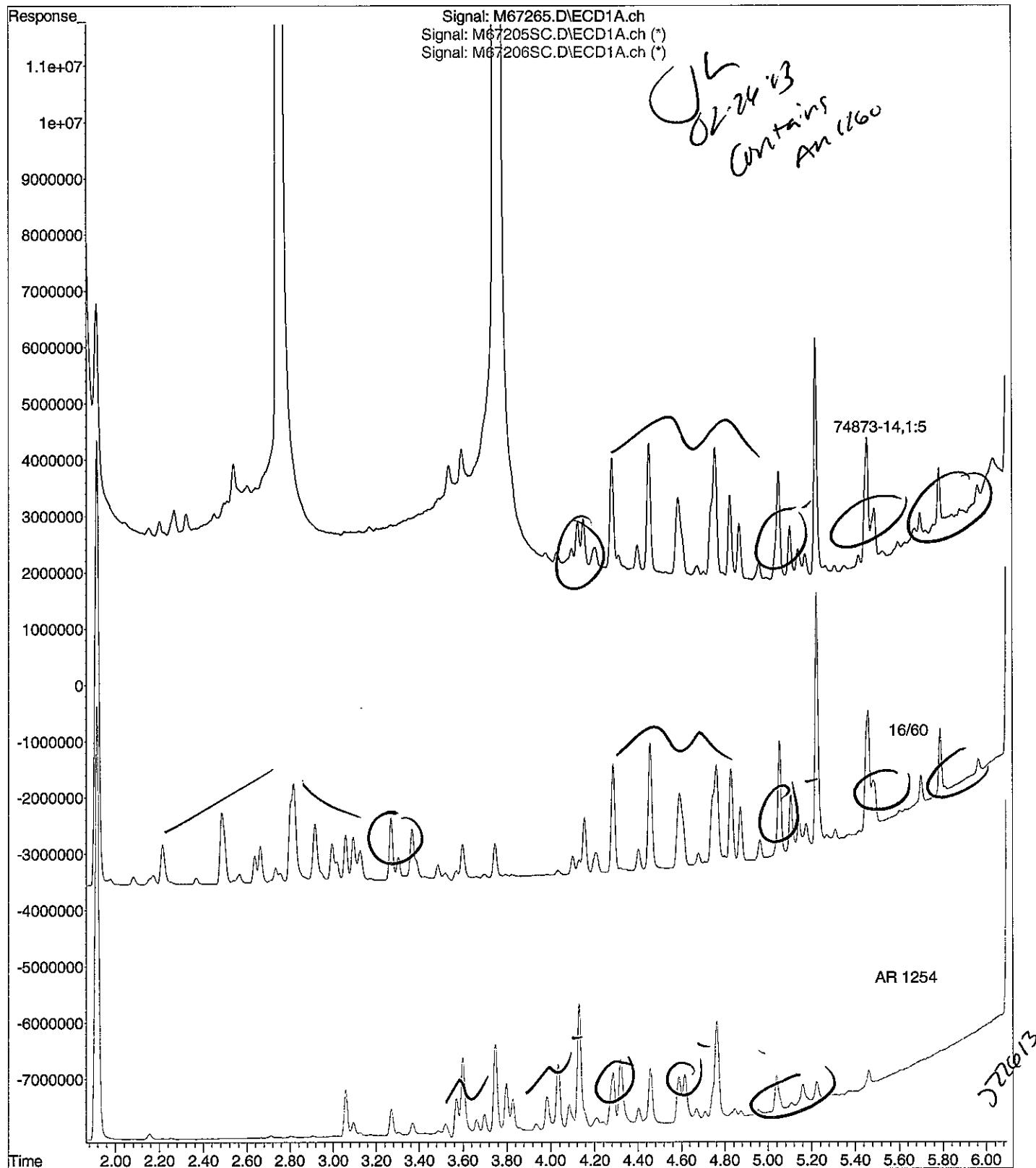
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67265.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 9:10 pm
 Operator : JK
 Sample : 74873-14,1:5,,A/C
 Misc : SOIL
 ALS Vial : 18 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 12:51:55 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:29 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67265.D
Operator : JK
Acquired : 22 Feb 2013 9:10 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-14,1:5,,A/C
Misc Info : SOIL
Vial Number: 18



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February 26, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-213 (4-6')

Lab Sample ID: 74873-15
Matrix: Solid
Percent Solid: 91
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	1790

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	103	%
Decachlorobiphenyl	59	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-15,1:5,,A/C

Column ID: 0.25 mm

Data File: M67266.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1791	1562	13.7	

Column to be used to flag RPD values greater than QC limit of 40%

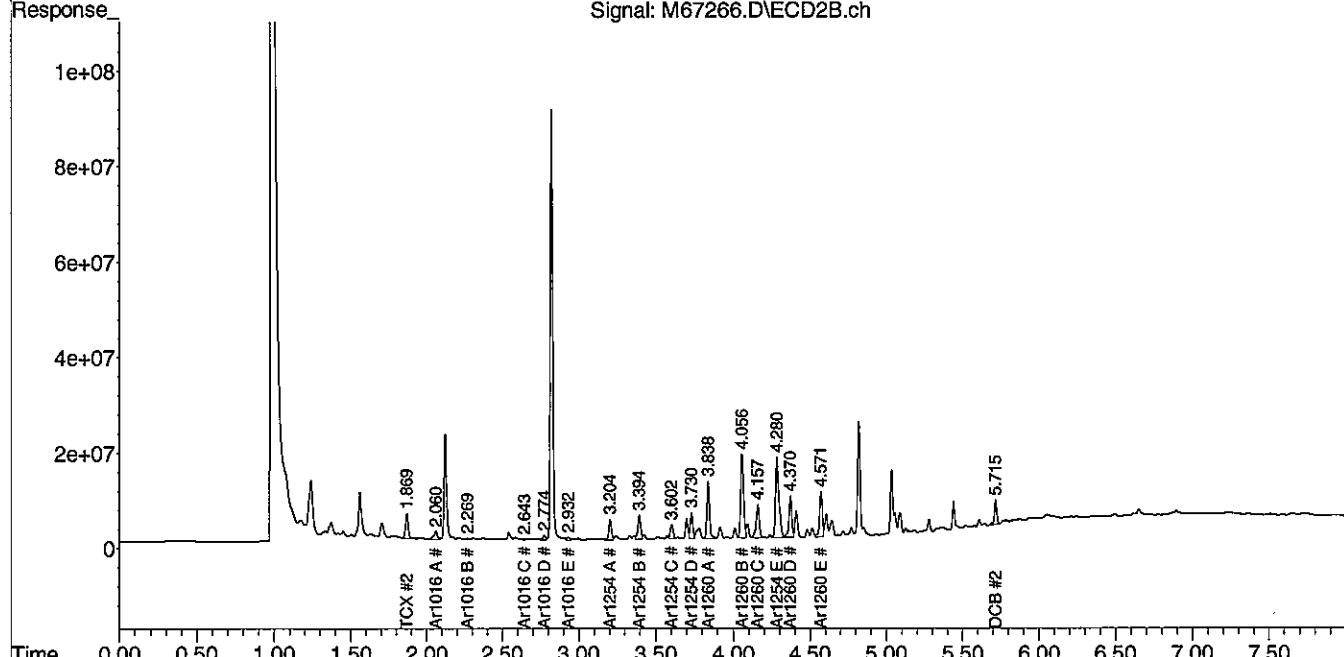
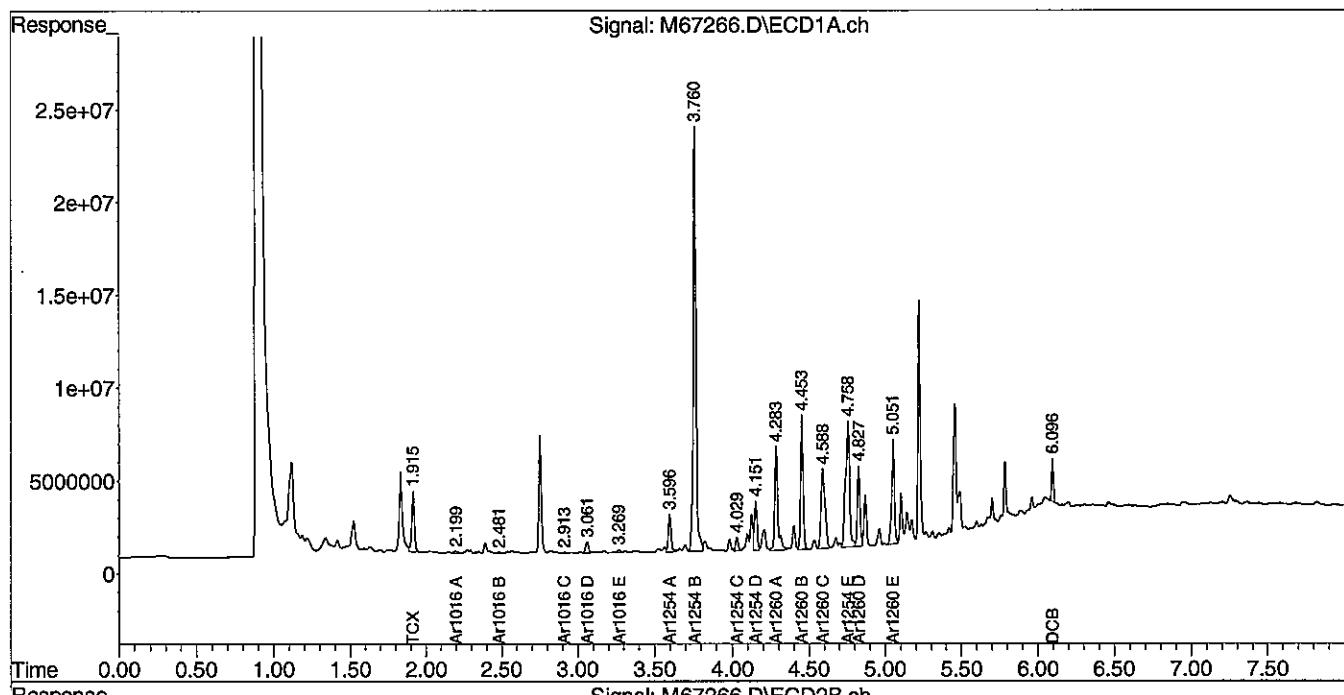
* Values outside QC limits

Comments: _____

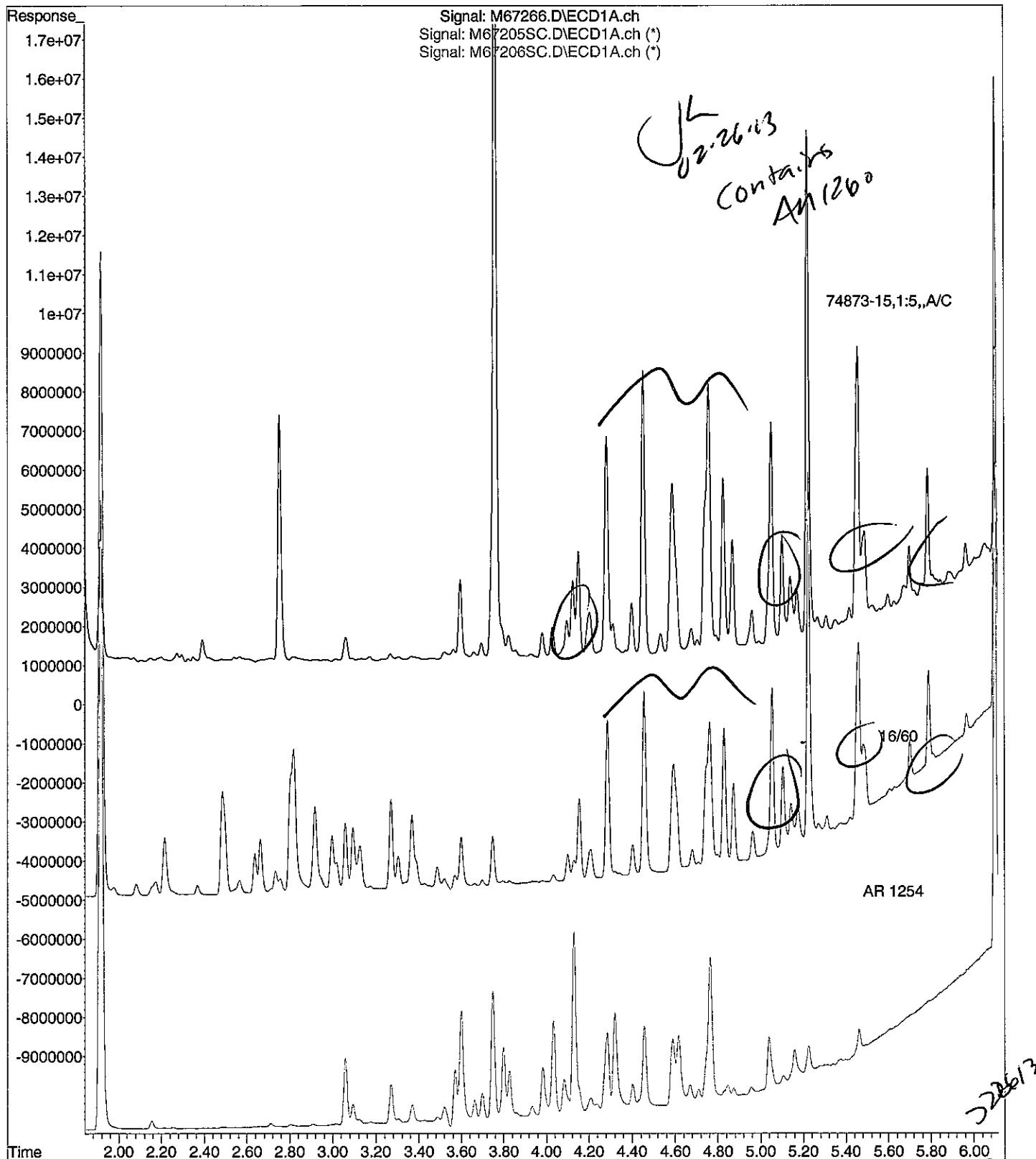
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67266.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 9:20 pm
 Operator : JK
 Sample : 74873-15,1:5,,A/C
 Misc : SOIL
 ALS Vial : 19 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 12:51:21 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67266.D
Operator : JK
Acquired : 22 Feb 2013 9:20 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-15,1:5,,A/C
Misc Info : SOIL
Vial Number: 19



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-213 (6-7.5')

Lab Sample ID:	74873-16
Matrix:	Solid
Percent Solid:	90
Dilution Factor:	11
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	9140

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	99	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-16,1:10,,A/C

Column ID: 0.25 mm

Data File: M67423.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.7

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	9144	7785	16.0	

Column to be used to flag RPD values greater than QC limit of 40%

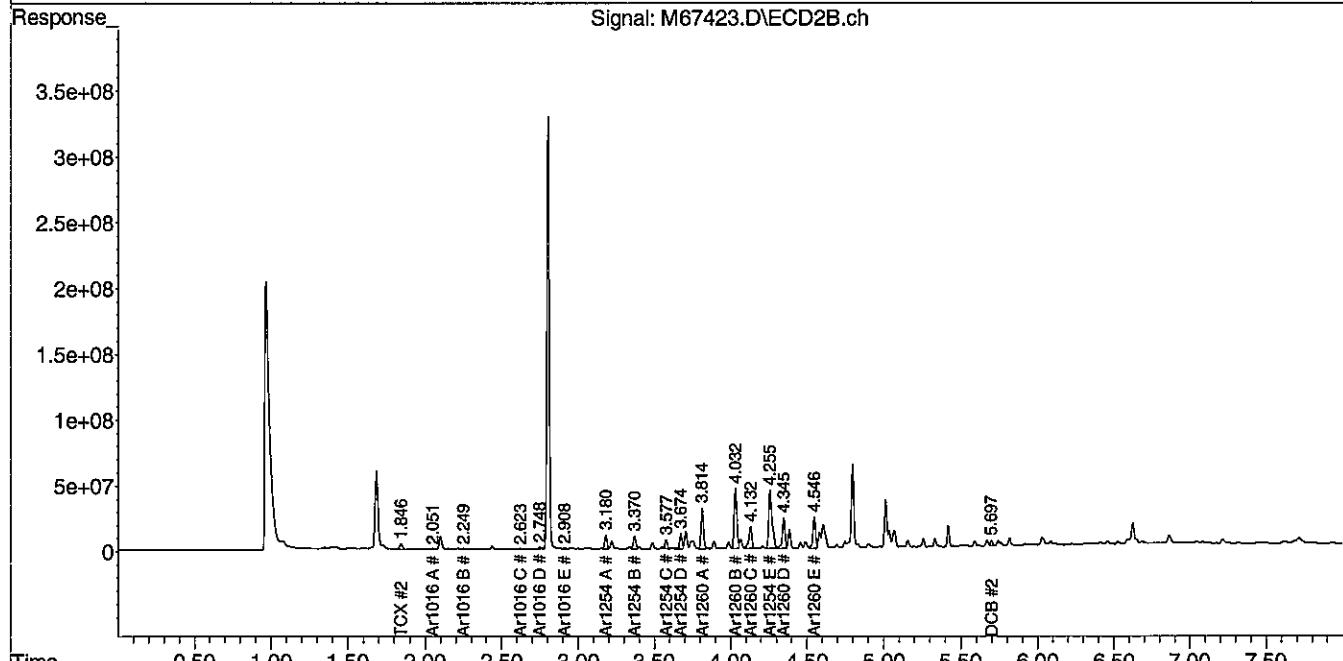
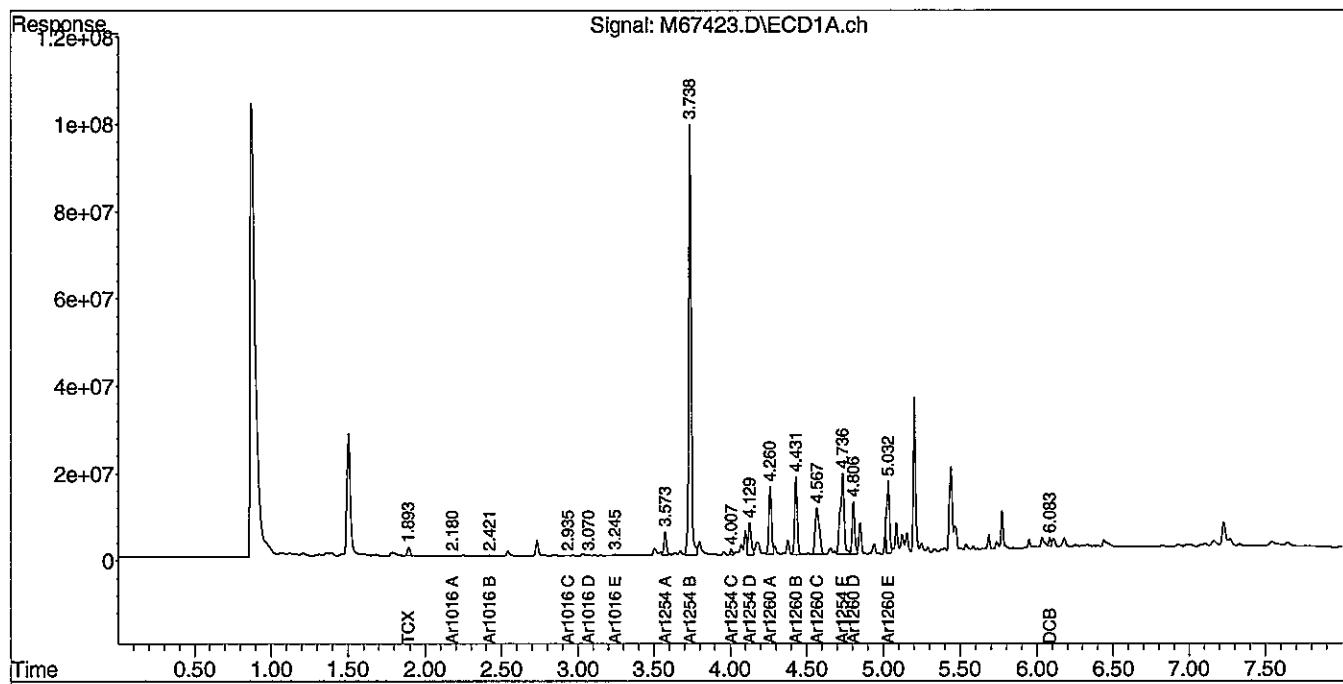
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67423.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 8:34 pm
 Operator : JK
 Sample : 74873-16,1:10,,A/C
 Misc : SOIL
 ALS Vial : 15 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:49:10 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 26, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-217 (0-2")

Lab Sample ID:	74873-17
Matrix:	Solid
Percent Solid:	53
Dilution Factor:	9
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	297	U
PCB-1221	297	U
PCB-1232	297	U
PCB-1242	297	U
PCB-1248	297	U
PCB-1254	297	U
PCB-1260	297	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	87	%
Decachlorobiphenyl	80	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

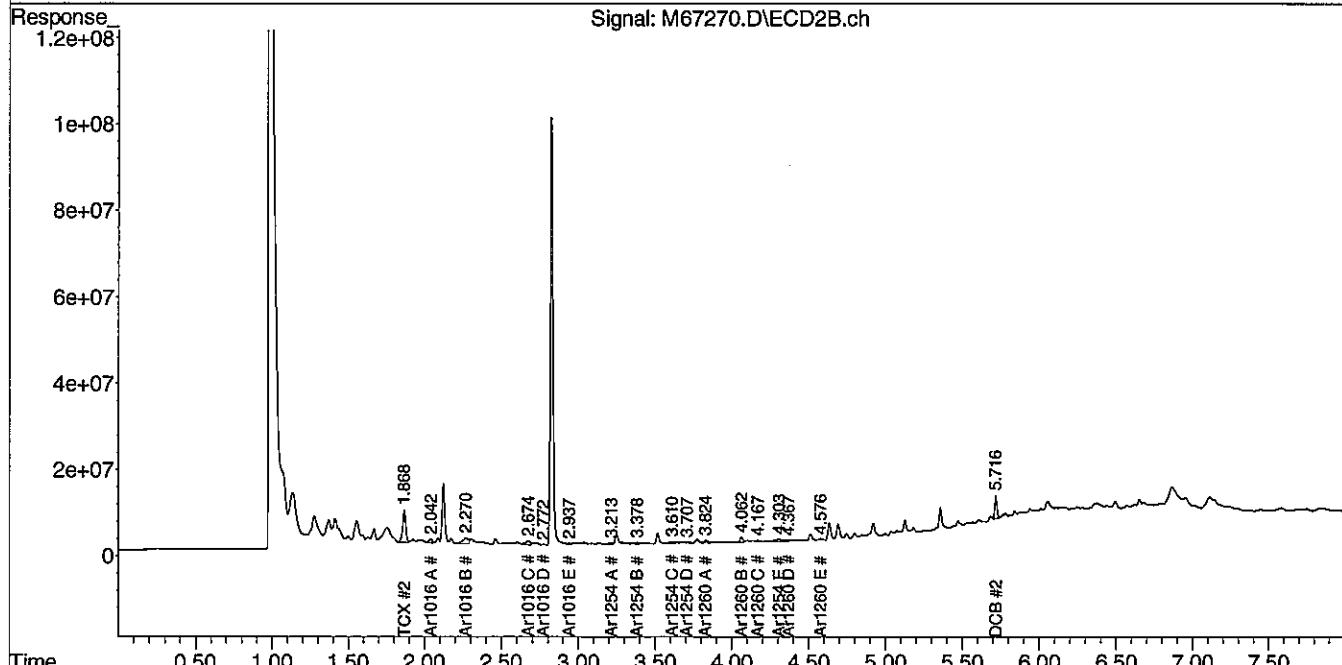
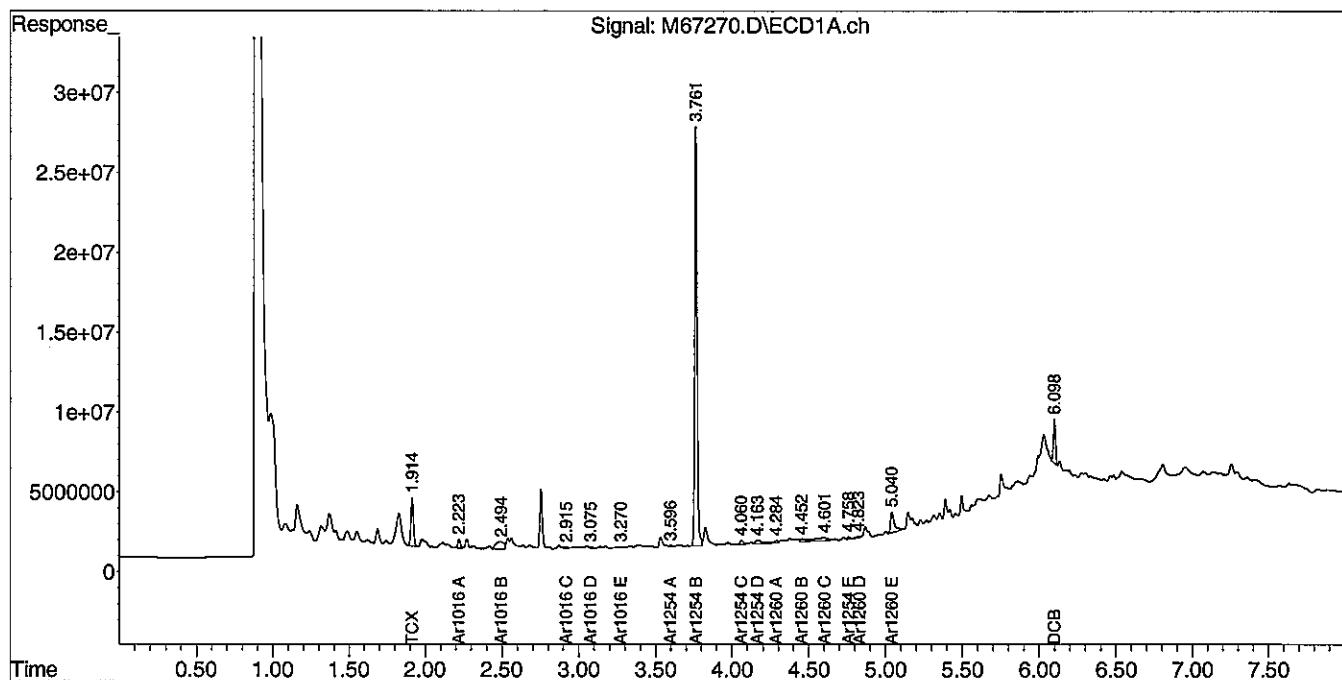
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * Quantitation limits increased due to the sample matrix affect.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67270.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 10:00 pm
 Operator : JK
 Sample : 74873-17,1:5,,A/C
 Misc : SOIL
 ALS Vial : 23 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 12:57:42 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID		Lab Sample ID:	74873-18
Project Name:	MAINE ENERGY	Matrix:	Solid
Project Number:	12-3259.1	Percent Solid:	86
Field Sample ID:	SE-SB-217 (2-4")	Dilution Factor:	6
		Collection Date:	02/19/13
		Lab Receipt Date:	02/20/13
		Extraction Date:	02/22/13
		Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	1160

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	134*	%
Decachlorobiphenyl	94	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * Surrogate recovery outside control limits. Secondary surrogate is in control.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-18,1:5,,A/C

Column ID: 0.25 mm

Data File: M67457.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1159	943	20.6	

Column to be used to flag RPD values greater than QC limit of 40%

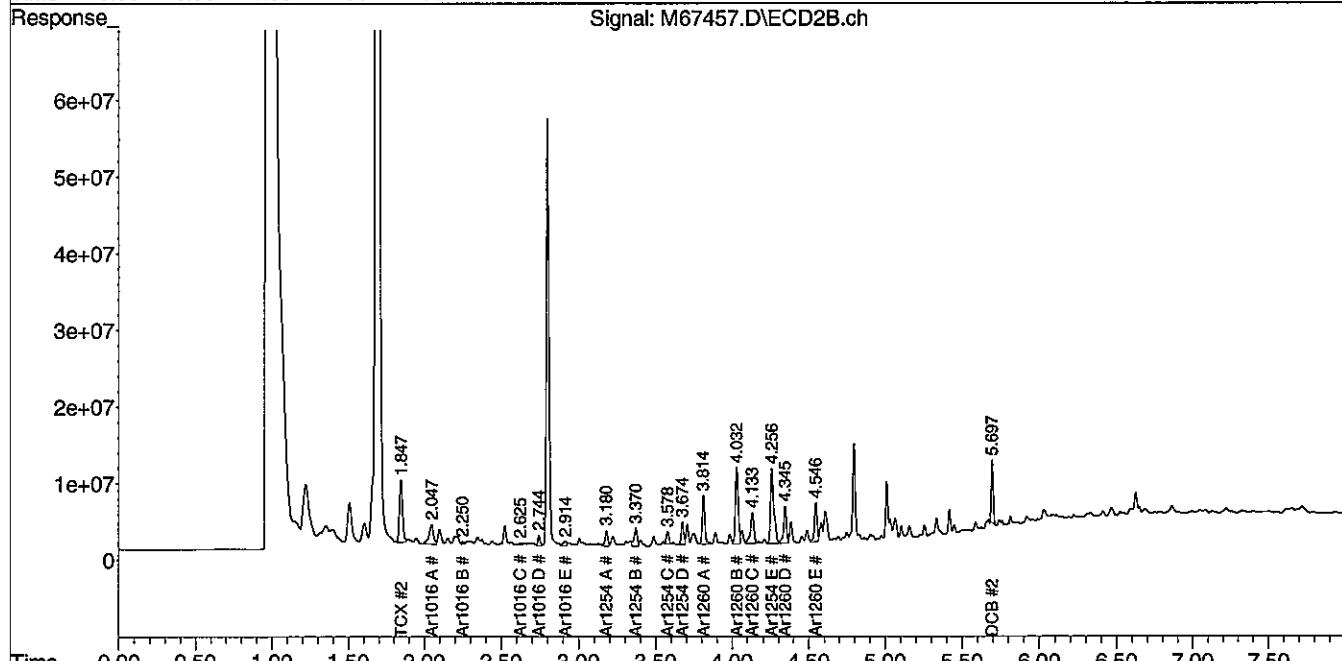
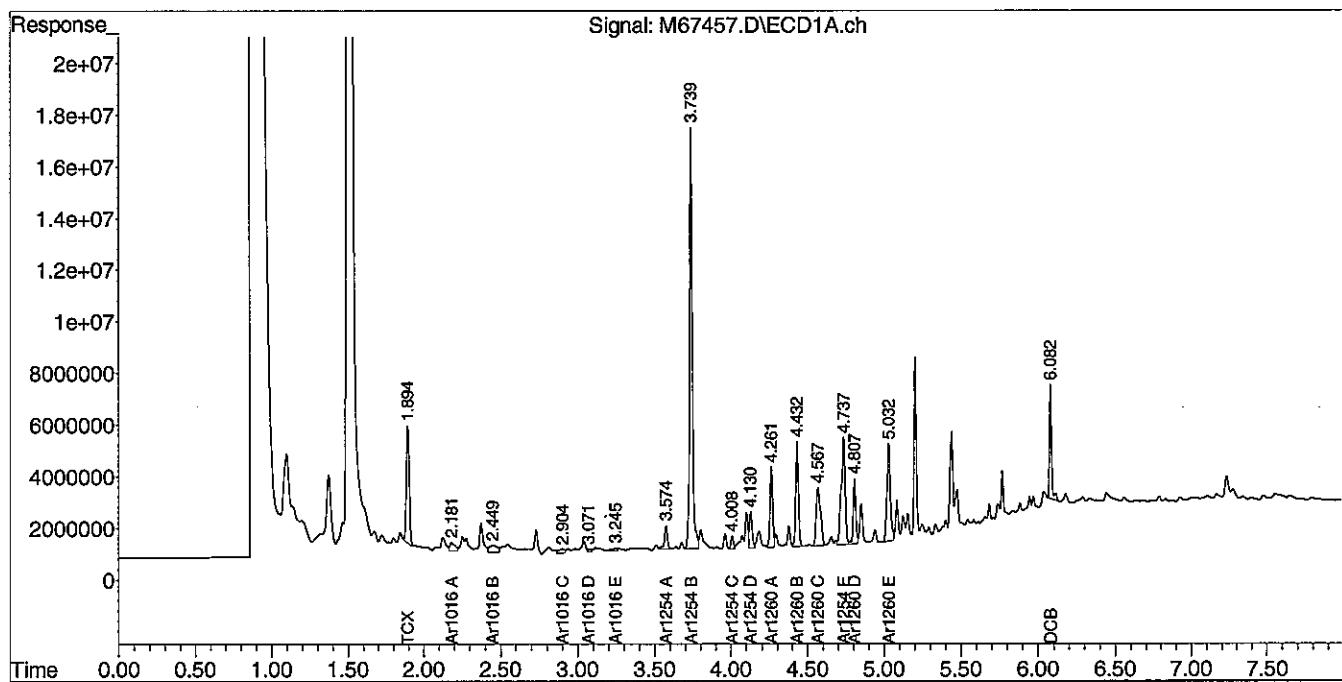
* Values outside QC limits

Comments: _____

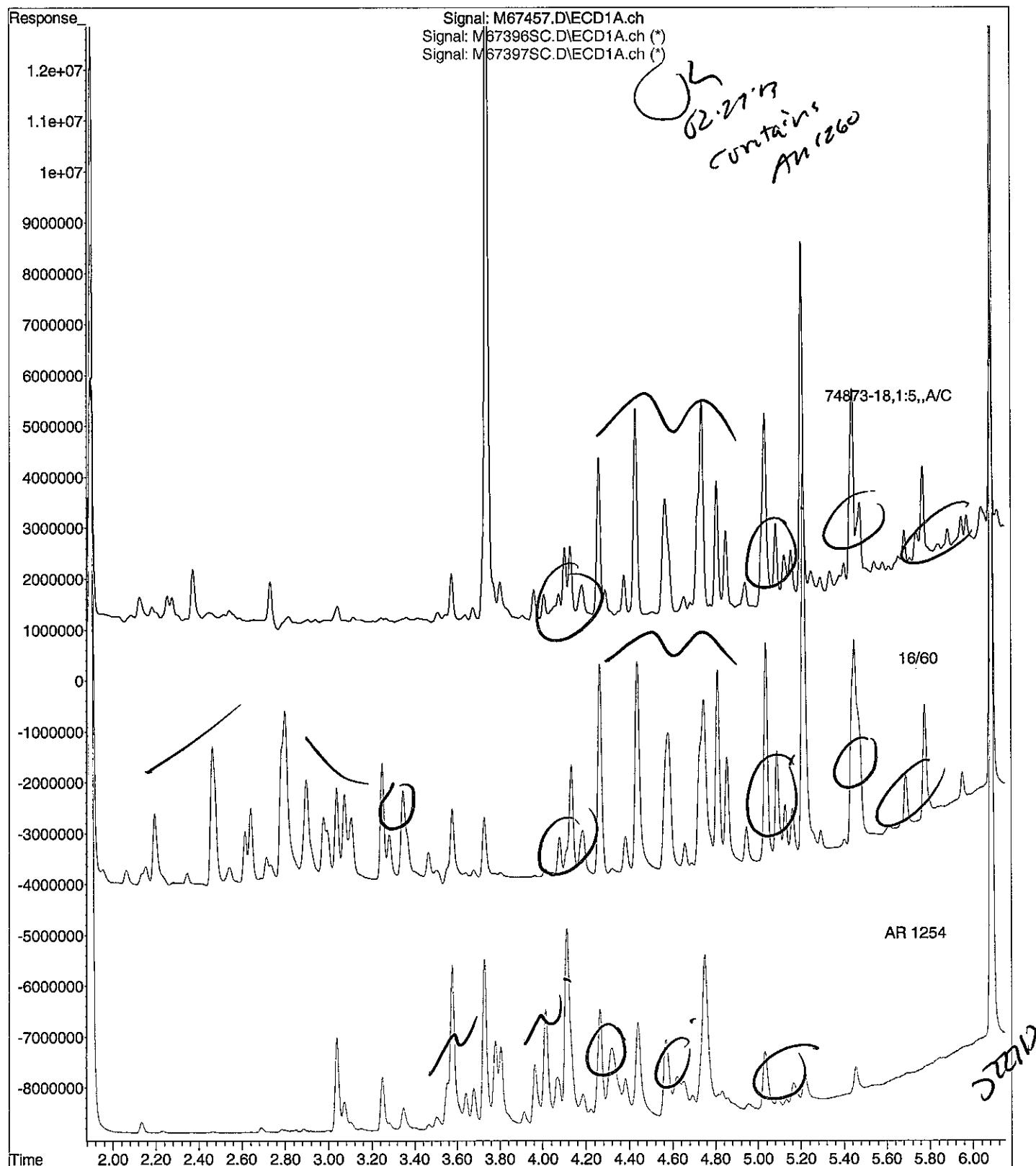
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67457.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 2:16 am
 Operator : JK
 Sample : 74873-18,1:5,,A/C
 Misc : SOIL
 ALS Vial : 41 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:49:43 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67457.D
Operator : JK
Acquired : 27 Feb 2013 2:16 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-18,1:5,,A/C
Misc Info : SOIL
Vial Number: 41



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-217 (4-6')

Lab Sample ID: 74873-19
Matrix: Solid
Percent Solid: 93
Dilution Factor: 1.0
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	853

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	80	%
Decachlorobiphenyl	72	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-19,,A/C

Column ID: 0.25 mm

Data File: M67458.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	853	726	16.1	

Column to be used to flag RPD values greater than QC limit of 40%

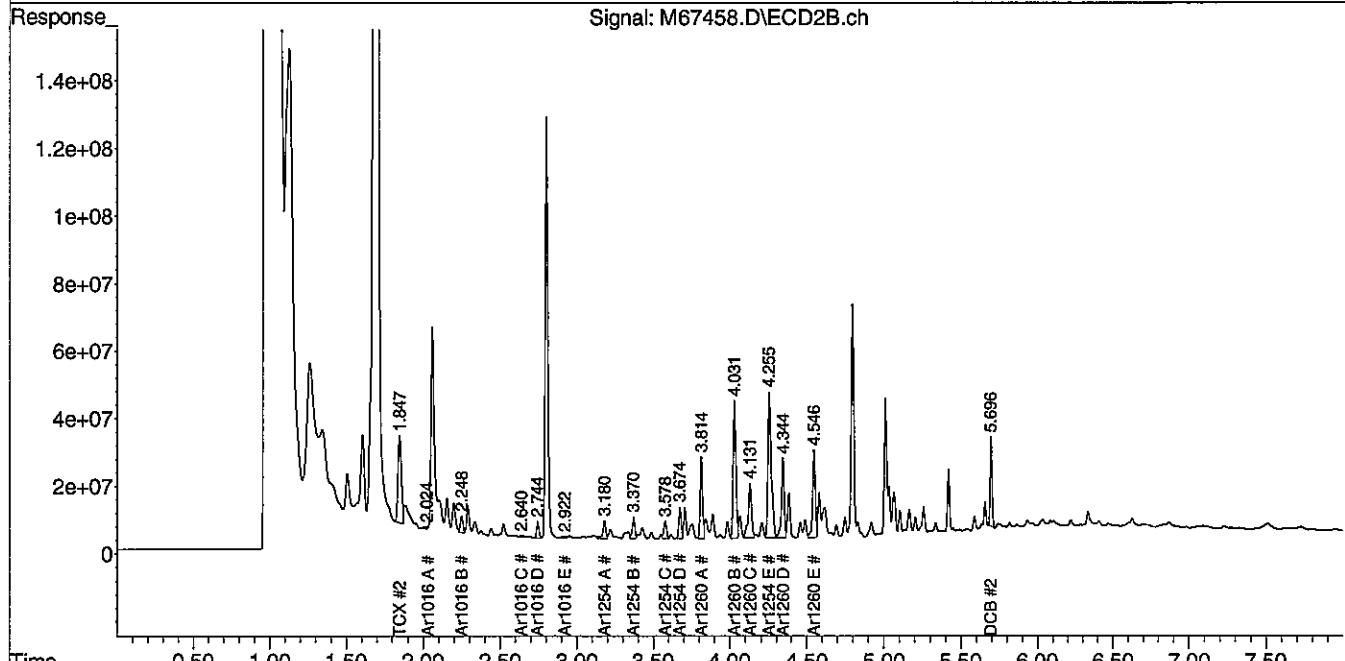
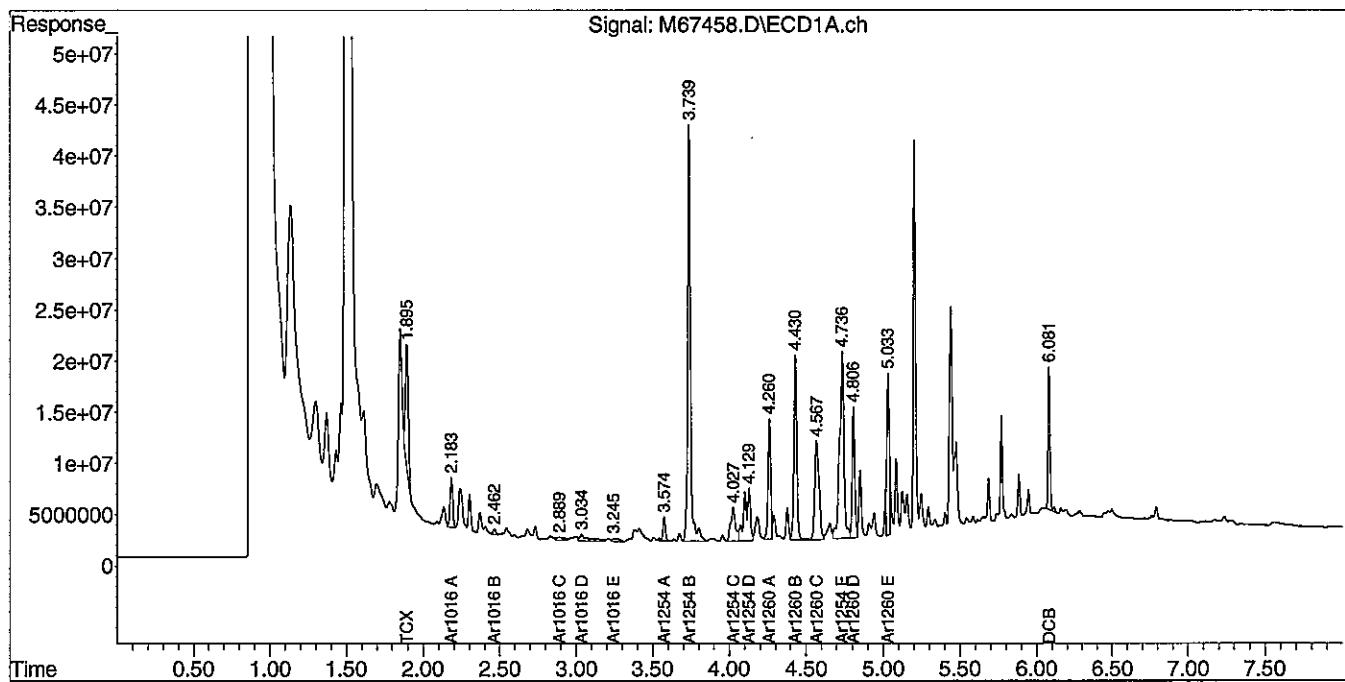
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67458.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 2:26 am
 Operator : JK
 Sample : 74873-19,,A/C
 Misc : SOIL
 ALS Vial : 42 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:51:35 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-217 (8-10')

Lab Sample ID: 74873-20
Matrix: Solid
Percent Solid: 89
Dilution Factor: 11
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	7860
PCB-1260	363	6680

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	115	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74873

GC Column #1: STX-CLPesticides I

Sample: 74873-20,1:10,,A/C

Column ID: 0.25 mm

Data File: M67504.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.7

Column ID: 0.25 mm

COMPOUND	Column #1		Column #2	
	SAMPLE RESULT (ug/kg)		SAMPLE RESULT (ug/kg)	RPD #
PCB 1260	6684		5068	27.5
PCB 1254	7859		6730	15.5

Column to be used to flag RPD values greater than QC limit of 40%

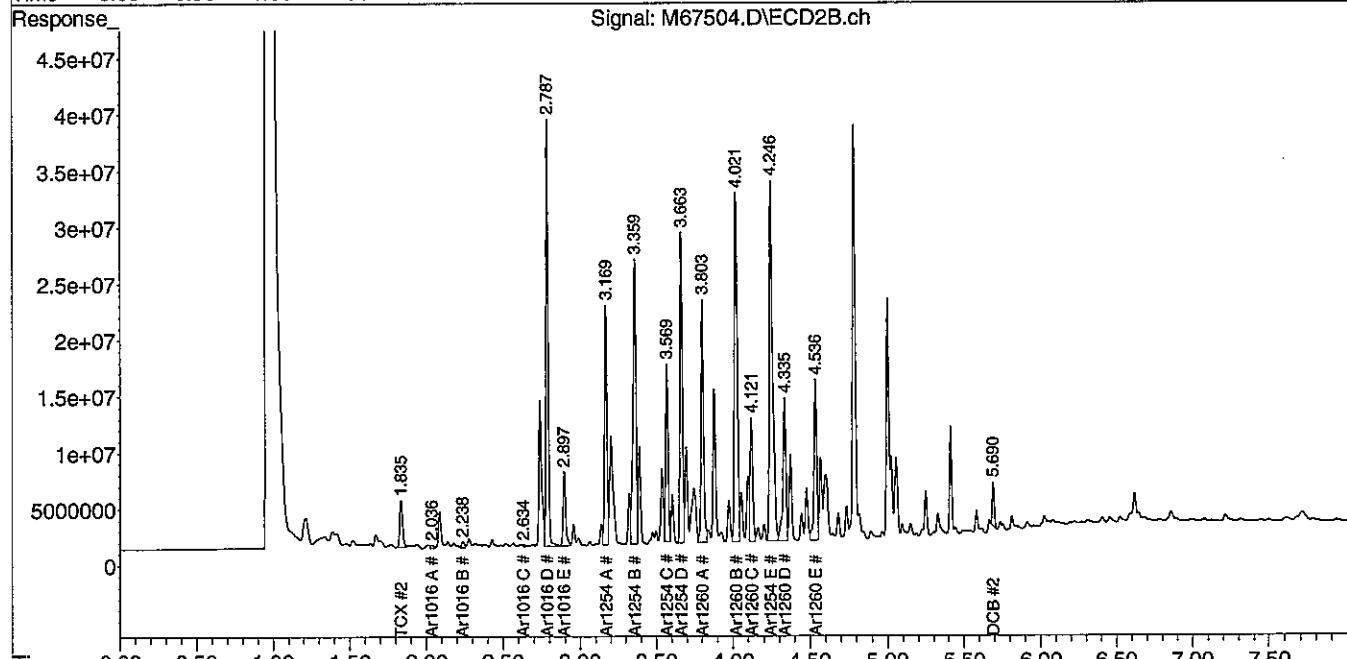
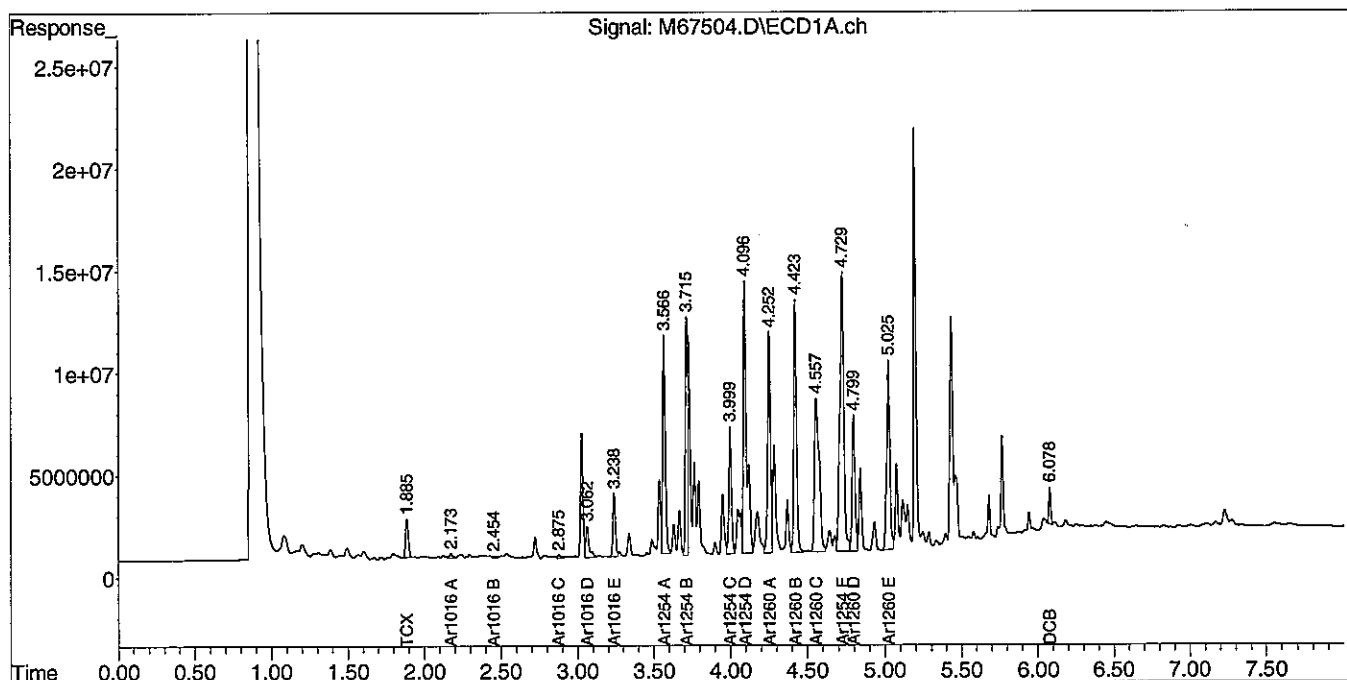
* Values outside QC limits

Comments: _____

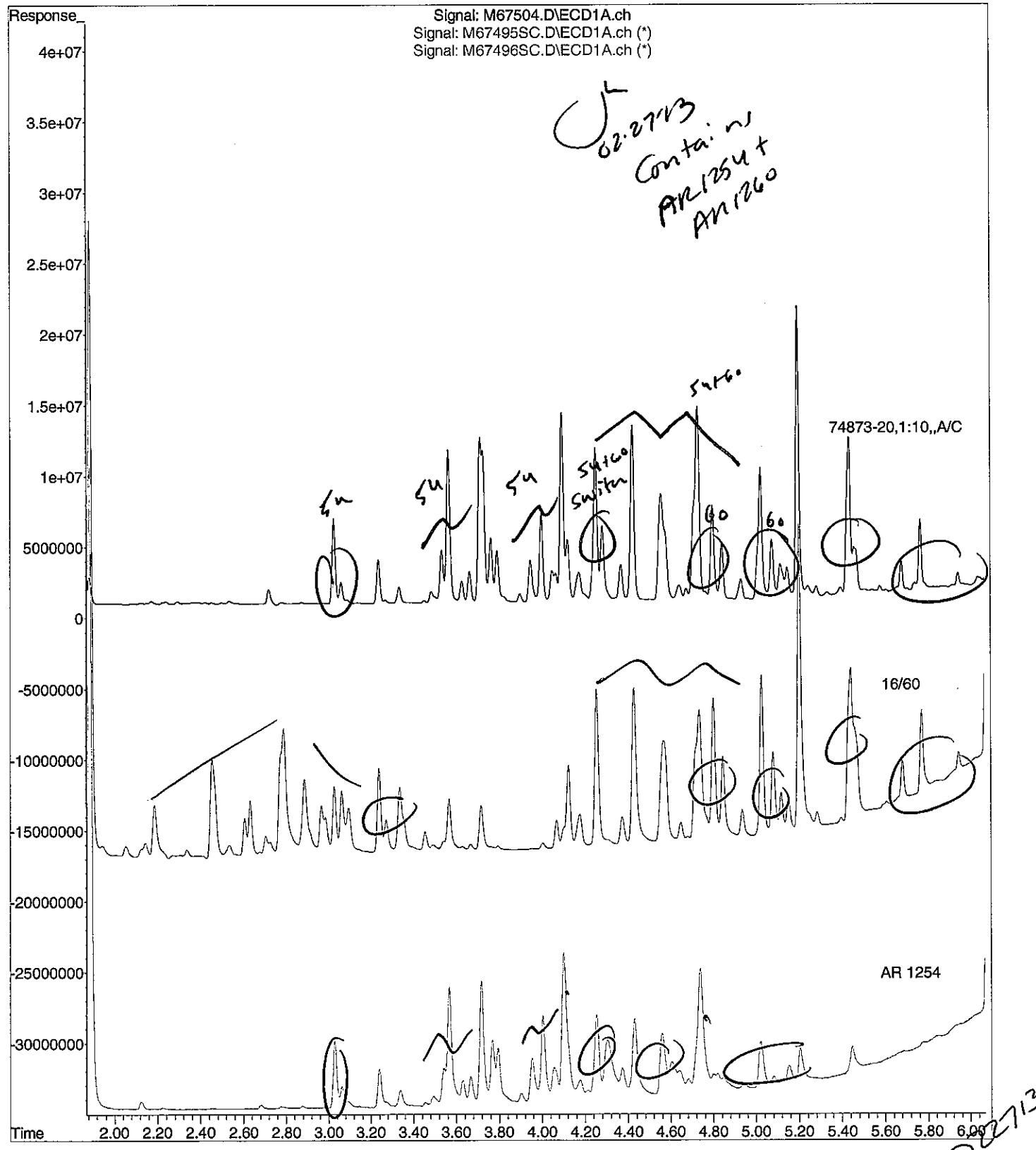
Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67504.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:29 pm
 Operator : JK
 Sample : 74873-20,1:10,,A/C
 Misc : SOIL
 ALS Vial : 7 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 16:07:39 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022713-M\M67504.D
Operator : JK
Acquired : 27 Feb 2013 1:29 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74873-20,1:10,,A/C
Misc Info : SOIL
Vial Number: 7



PCB
QC FORMS

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February 22, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	Lab QC

Lab Sample ID:	B022013PSOX2 RR
Matrix:	Soil
Percent Solid:	100
Dilution Factor:	1.0
Collection Date:	
Lab Receipt Date:	
Extraction Date:	02/20/13
Analysis Date:	02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	103	%
Decachlorobiphenyl	94	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

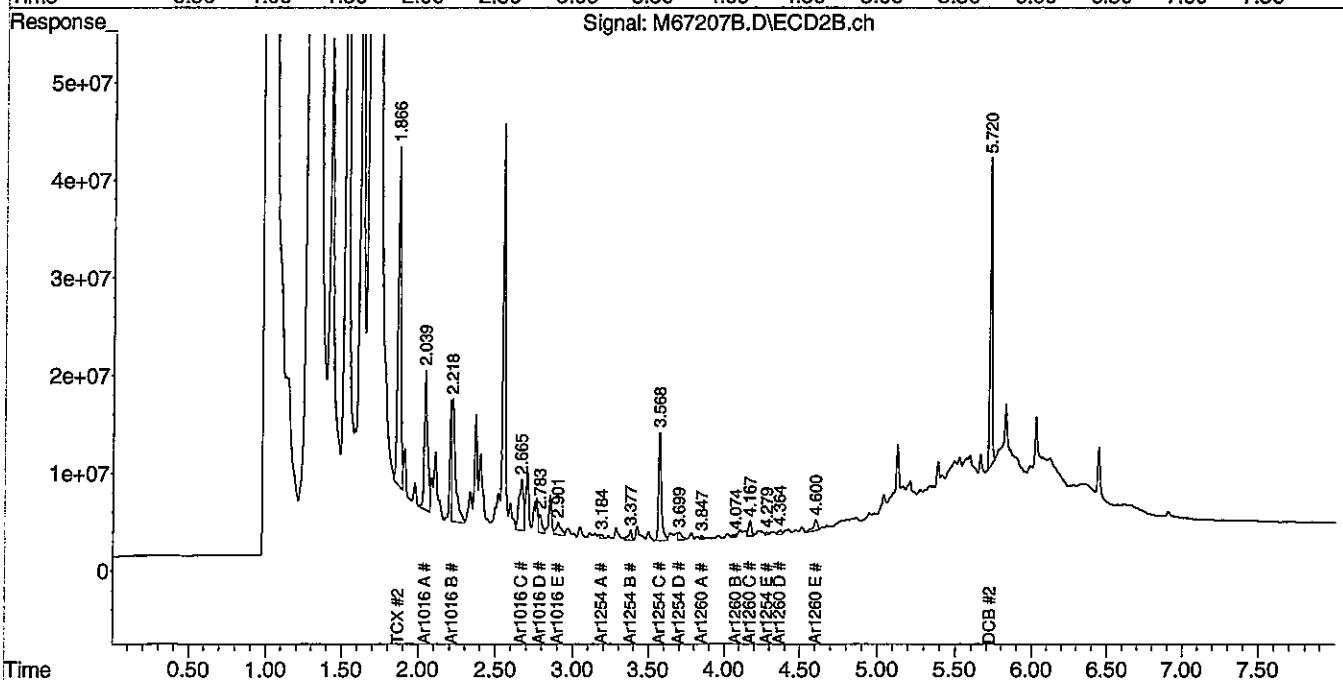
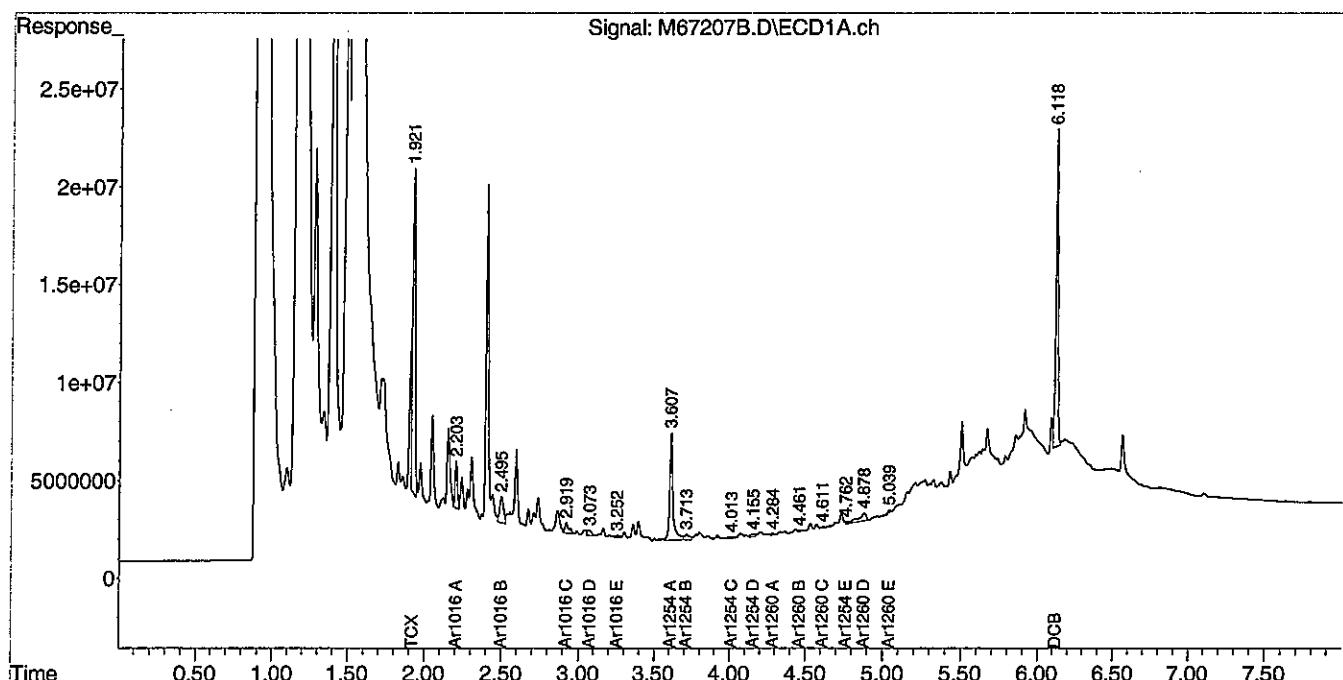
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67207B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 10:06 am
 Operator : JK
 Sample : B022013PSOX2,RR,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 22 15:58:42 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	Lab QC

Lab Sample ID:	B022513PSOX
Matrix:	Soil
Percent Solid:	100
Dilution Factor:	1.0
Collection Date:	
Lab Receipt Date:	
Extraction Date:	02/25/13
Analysis Date:	02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	182*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

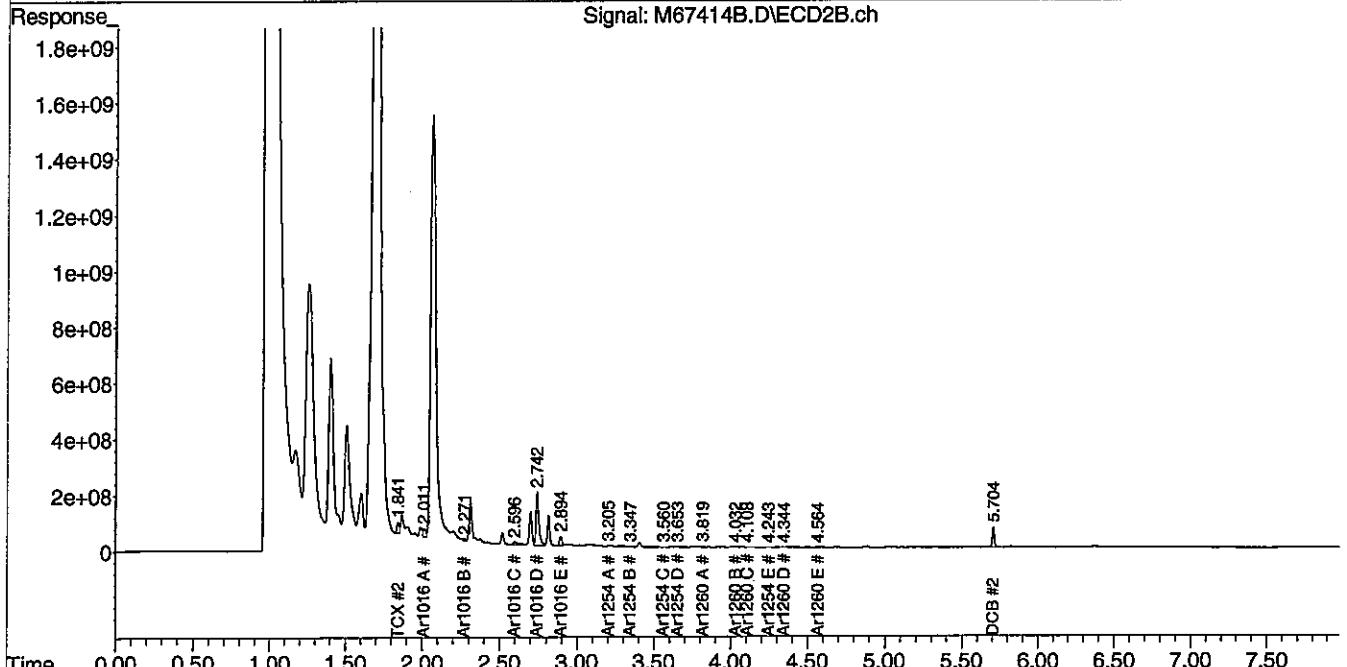
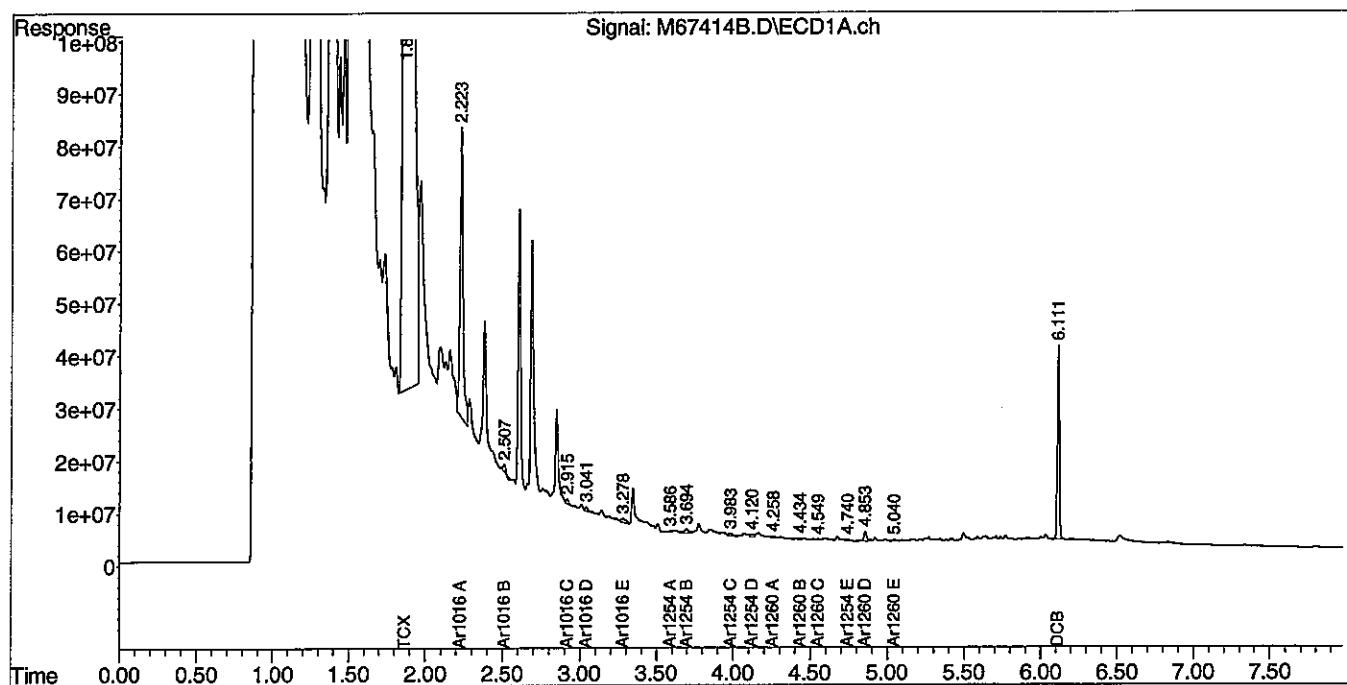
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * Surrogate recovery outside control limits. Secondary surrogate is in control.

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67414B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 7:03 pm
 Operator : JK
 Sample : B022513PSOX,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 10:22:59 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 26, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	Lab QC

Lab Sample ID:	B022113PSOX2
Matrix:	Soil
Percent Solid:	100
Dilution Factor:	1.0
Collection Date:	
Lab Receipt Date:	
Extraction Date:	02/21/13
Analysis Date:	02/22/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	70	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

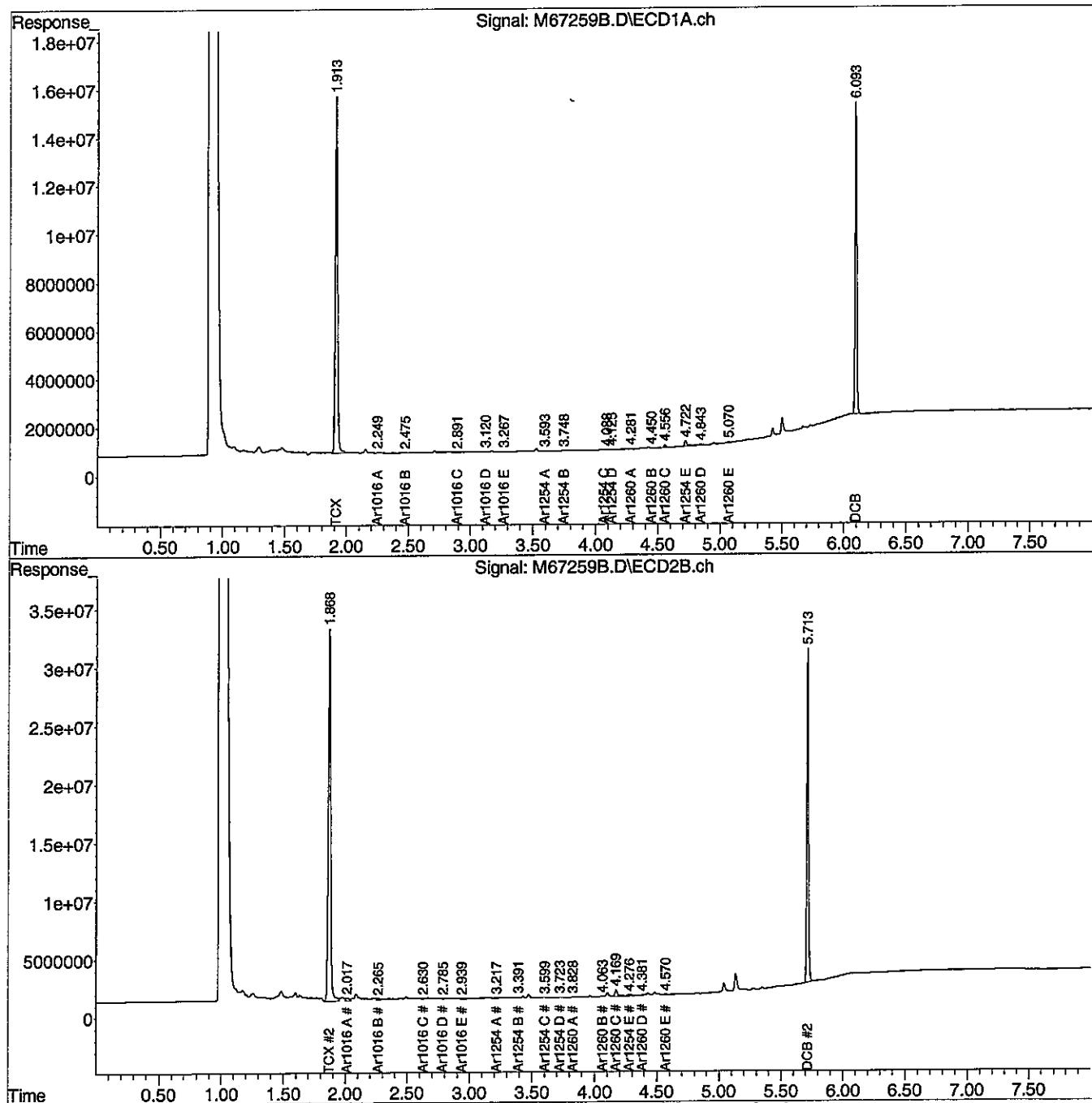
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67259B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 22 Feb 2013 8:09 pm
 Operator : JK
 Sample : B022113PSOX2,,A/C
 Misc : SOIL
 ALS Vial : 12 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 11:43:01 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022213PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	69	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

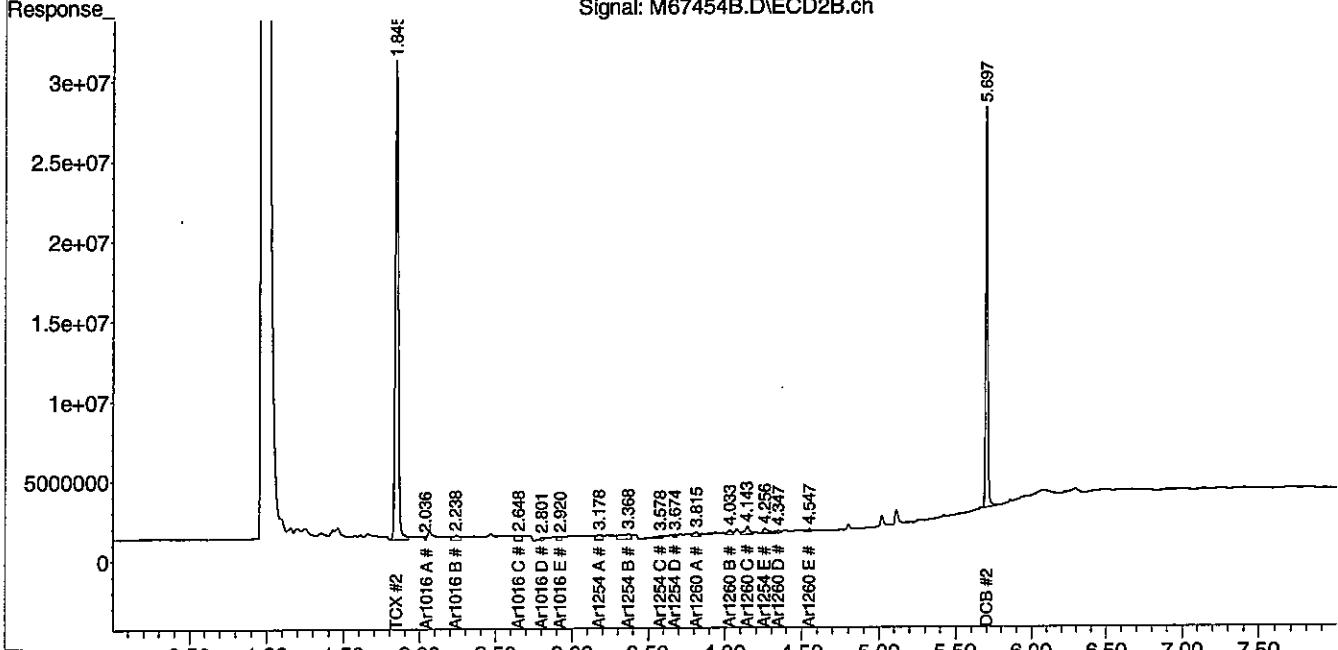
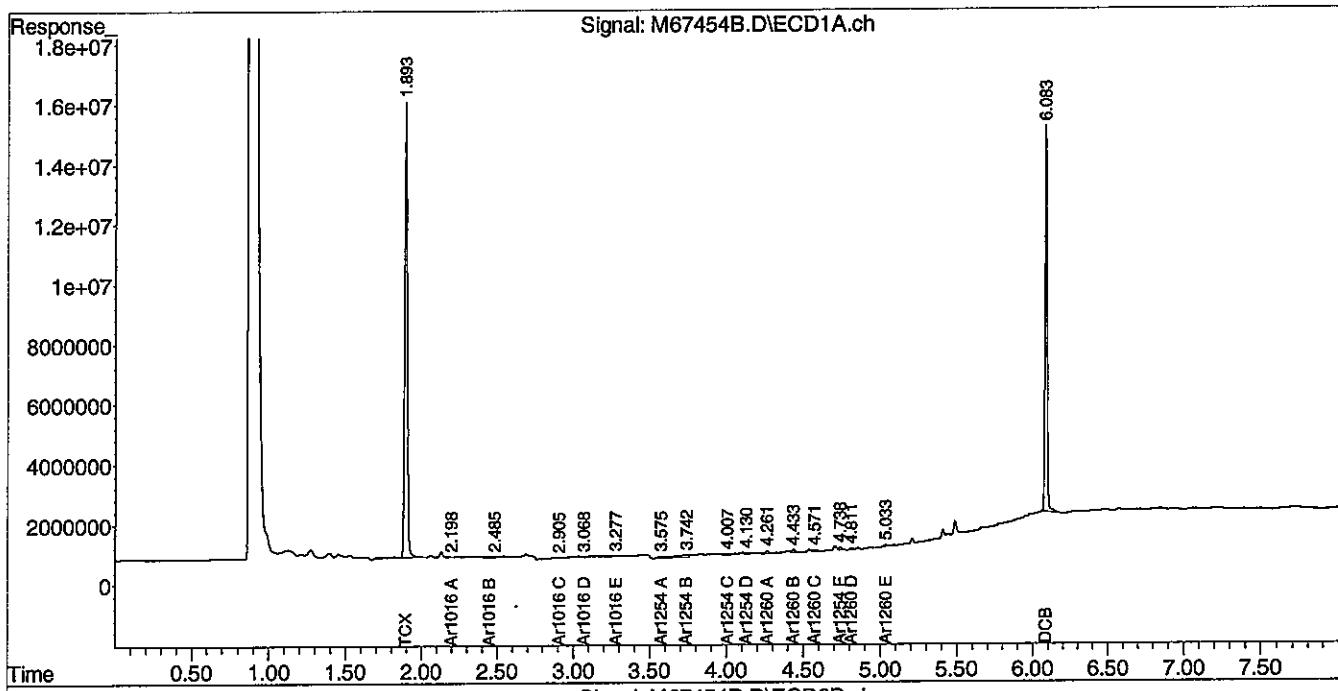
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67454B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:46 am *Feb 27 2013*
 Operator : JK
 Sample : B022213PSOX,, A/C (Sig #1); B022113PSOX,, A/C (Sig #2)
 Misc : SOIL
 ALS Vial : 38 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:47:07 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	Lab QC

Lab Sample ID:	B022213PSOX RR
Matrix:	Soil
Percent Solid:	100
Dilution Factor:	1.0
Collection Date:	
Lab Receipt Date:	
Extraction Date:	02/21/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	75	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

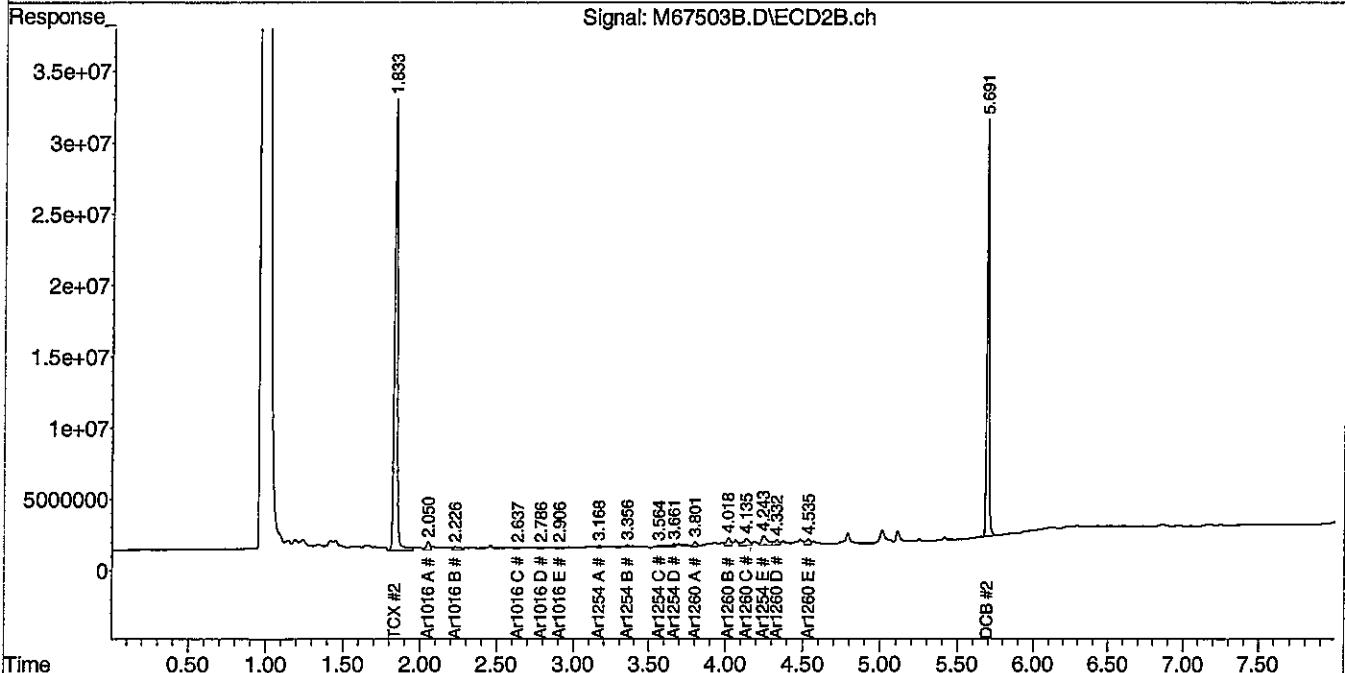
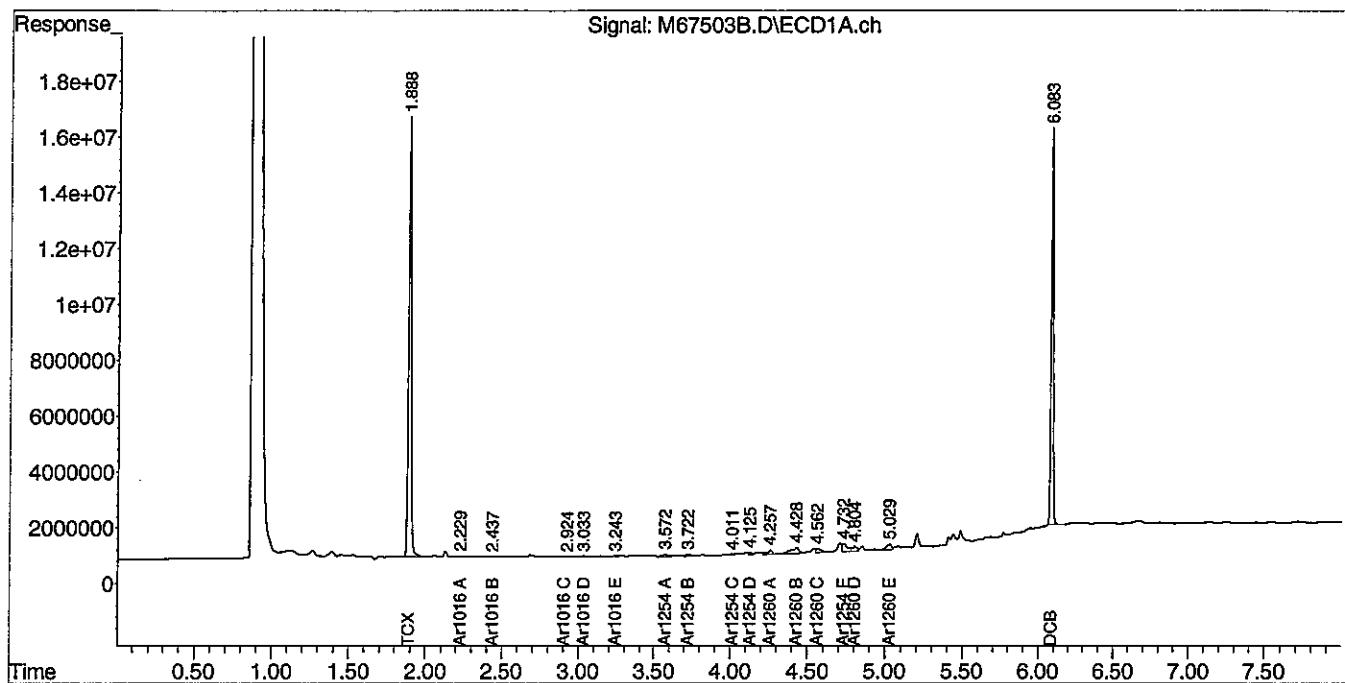
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67503B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:19 pm
 Operator : JK
 Sample : B022213PSOX,RR2,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:19 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74873

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits
* Values outside QC limits
D System Monitoring Compound diluted out

PCB SOIL SYSTEM MONITORING COMPOUNDS SUMMARY

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74873

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D. System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74873

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74873

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits
* Values outside QC limits
D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74873

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits
* Values outside QC limits
D System Monitoring Compound diluted out

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022013PSOX2,RR,,A/C

Spike: L022013PSOX2,RR,,A/C

Spike duplicate: LD022013PSOX2,RR,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	157	78		155	78		0.8	
PCB 1260	200	200	60	130	30	0	184	92		149	74		21.3	
PCB 1016 #2	200	200	65	140	30	0	163	81		165	83		1.4	
PCB 1260 #2	200	200	60	130	30	0	194	97		147	74		27.5	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

**PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: 74873-1,RR,,A/C

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: 74873-1,MS,RR,,A/C

Spike duplicate: 74873-1,MSD,RR,,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP	RPD	#
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#
PCB 1016	209	202	60	140	30	0	180	86		170	84	5.6
PCB 1260	209	202	60	140	30	65	241	84		338	135	33.6 *
PCB 1016 #2	209	202	60	140	30	0	153	73		253	125	49.2 *
PCB 1260 #2	209	202	60	140	30	64	333	129		252	93	27.7

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG:

Column ID: 0.25 mm

Non-spiked sample: B022113PSOX2,,A/C

GC Column #2: STX-CLPesticides II

Spike: L022113PSOX2,,A/C

Column ID: 0.25 mm

Spike duplicate: LD022113PSOX2,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	169	84		166	83		1.5	
PCB 1260	200	200	60	130	30	0	162	81		160	80		1.5	
PCB 1016 #2	200	200	65	140	30	0	166	83		164	82		1.0	
PCB 1260 #2	200	200	60	130	30	0	160	80		161	80		0.4	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022513PSOX,,A/C

Spike: L022513PSOX,,A/C

Spike duplicate: LD022513PSOX,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	179	90		192	96		7.0	
PCB 1260	200	200	60	130	30	0	207	103		200	100		3.6	
PCB 1016 #2	200	200	65	140	30	0	171	86		169	85		1.0	
PCB 1260 #2	200	200	60	130	30	0	197	99		186	93		5.9	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

**PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: 74873-16,1:10,,A/C

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: 74873-16,MS,1:100,,A/C

Spike duplicate: 74873-16,MSD,1:100,,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC #	RESULT (ug/kg)	% REC #	RPD #
PCB 1016	221	210	65	140	30	0	D	D	D	D	
PCB 1260	221	210	60	130	30	9144	68555	26827 *	8979	-79 *	153.7 *
PCB 1016 #2	221	210	65	140	30	0	D	D	D	D	
PCB 1260 #2	221	210	60	130	30	7785	57720	22547 *	7607	-85 *	153.4 *

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022213PSOX,,A/C

Spike: L022213PSOX,,A/C

Spike duplicate: LD022213PSOX,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	178	89		175	87		2.0	
PCB 1260	200	200	60	130	30	0	179	90		173	87		3.3	
PCB 1016 #2	200	200	65	140	30	0	173	86		166	83		4.0	
PCB 1260 #2	200	200	60	130	30	0	172	86		166	83		3.7	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form

CINCHYDICS environmental laboratory LLC

Environmental
laboratory LLC

 Project Name: MAINE ENERGY Project #: 12-3259.1		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax	
Samples were: 1) Shipped or hand-delivered <input checked="" type="checkbox"/> 2) Temperature (°C): <u>03-07.0</u> 3) Received in good condition: <input checked="" type="checkbox"/> N/A 4) pH checked by: <u>SP2001/3</u> 5) Labels checked by:					
Matrix Key: C = Concrete WP = Wipe DW = Drinking Water S = Soil / Sludge SW = Surface Water O = Oil X = Other E = Extract					
Circle and/or Write Required Analysis Followed by Preservation Code Please fill in preservation code here					
Sample Identification	Sample Date	Sample Time	VOC: 8260 524.2 624	SVOCl: 8270 625 PAH only SIM	Pesticides: 8081 608
SE-SB-216 (2.5"-24")	2-19	1058	N	N	X
SE-SB-216 (8-4")		1120	Z	X	X
SE-SB-216 (4-6")		1125	N	X	X
SE-SB-215 (0-2")		1136	N	X	X
SE-SB-215 (3-4")		1145	N	X	X
SE-SB-215 (4-6")		1154	N	X	X
SE-SB-215 (6-8")		1205	N	X	X
SE-SB-215 (8-16")		1265	N	X	X
SE-SB-214 (0-2")		1230	Z	X	X
SE-SB-214 (2-4")		1238	N	X	X
** List requested metals here Project Requirements: *Fee may apply					
Comments, Additional Analyses, or Special Instructions: <u>Comments: No comments.</u>					
Email Results to: <u>JLAW@GEOGRAPHICS.NH.COM</u>					
Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> 4 Days* <input checked="" type="checkbox"/> 5 Days <input type="checkbox"/> Standard (8-10 business days)					
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.					
*Fee may apply; lab approval required					
Sampler Name (Print): <u>John Carson</u>					
Relinquished By Sampler: <u>TL</u>					
Relinquished By: _____					
Received By: <u>John Carson</u>					
Received By: _____					
Date: <u>2-20</u> Time: <u>1131</u> Received By: <u>John Carson</u>					
Date: _____ Time: _____ Received By: _____					
Date: _____ Time: _____ Received By: _____					
Report Type: <input type="checkbox"/> MCP* <input type="checkbox"/> Level II* <input checked="" type="checkbox"/> Standard <input type="checkbox"/> CTRCP* <input type="checkbox"/> Level III* <input type="checkbox"/> ME <input type="checkbox"/> DOD* <input type="checkbox"/> Level IV* <input type="checkbox"/> CT <input type="checkbox"/> Other: _____ <input type="checkbox"/> RI					
State Standard: <input type="checkbox"/> NH <input type="checkbox"/> MA <input checked="" type="checkbox"/> ME <input type="checkbox"/> (e.g. S-1 or GW-1) <input type="checkbox"/> EDD Required: <input checked="" type="checkbox"/> Y/N <input type="checkbox"/> Type: _____					

Analytics Report 74873 page 0096 of 98

Chain Of Custody Form

 environmental laboratory LLC		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax	For Analytics Use Only <i>03-07-05</i>	
Project Name:	MAINE ENERGY	Preservation Code:	Please <input checked="" type="checkbox"/> in preservation code here	Metals: RCRA P13 TAL23 Other: **	Samples were: 1) Shipped or hand-delivered 2) Temperature (°C): <i>15</i> 3) Received in good condition 4) pH checked by: <i>J. S. C.</i> 5) Labels checked by: <i>J. S. C.</i>	
Project#:	13-3259.1	Preservation Key: A = HCl B = 4°C C = Unpres D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other		VPH: Full or Ranges only TPH: 8015 (Diesel Range) ME217 TPH: 8015 (Gas Range) ME217 PCB: 8082 608 Sothlet V or N Pesticides: 8081 608 SVOC: 8270 625 PAH only SIM VOC: 8260 5242 624 Field Filtered Y or N	Matrix Key: C = Concrete W = Wipe WW = Wastewater SW = Surface Water E = Extract	
Company:	Sumit Corp.	Sample Date	Sample Time	Matrix	No. of Containers checked	
Report to:	John Crowley	SE-19-214 (4-6)	2-19 1234 N	S	1	
Address:	100 Main St. Lawrence, MA 01840	SE-19-214 (6-7)	1237 N	S	1	
Phone:	(201) 795-4882	SE-19-213 (0-2)	1247 N	S	1	
Quote #:		SE-19-213 (2-4)	1310 N	S	1	
PO# (if required):	12-3259.1	SE-19-213 (4-6)	1318 N	S	1	
		SE-19-213 (6-7.5)	1324 N	S	1	
		SE-19-217 (6-2)	1335 N	S	1	
		SE-19-217 (3-4)	1343 N	S	1	
		SE-19-217 (4-6)	1346 N	S	1	
		SE-19-217 (8-10)	1400 N	S	1	
** List requested metals here Project Requirements: *Fee may apply						
Email Results to:	<input type="checkbox"/> GinCityAnalytics.com <input type="checkbox"/> Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> 4 Days* <input type="checkbox"/> 5 Days <input checked="" type="checkbox"/> Standard <small>(*=10 business days)</small>					State Standard: <input type="checkbox"/> NH <input type="checkbox"/> MA <input checked="" type="checkbox"/> ME <small>(eg. S-1 or GW-1)</small> <input type="checkbox"/> CT <input type="checkbox"/> RI <input type="checkbox"/> Other: _____
Relinquished By Sampler:	<i>J. S. C.</i>					Report Type: <input type="checkbox"/> MCP* <input type="checkbox"/> Level II* <input type="checkbox"/> CTRCP* <input type="checkbox"/> Level III* <input type="checkbox"/> DOD* <input type="checkbox"/> Level IV* <input checked="" type="checkbox"/> Standard
Relinquished By:						Date: <i>2-20</i> Time: <i>1131</i> Received By: <i>K. Stinson</i>
Relinquished By:						Date: _____ Time: _____ Received By: _____



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 7478389 74873
CLIENT: Summit
PROJECT: Main Energy

COOLER NUMBER: 74,223
NUMBER OF COOLERS: 2

A: PRELIMINARY EXAMINATION:

1. Cooler received by(initials): EZ

DATE COOLER RECEIVED/OPENED: 02/20/13

2. Circle one:

Hand delivered
(If no, skip 3)

Shipped

3. Did cooler come with a shipping slip?

Y

N

3a. Enter carrier name and airbill number here: _____

4. Were custody seals on the outside of cooler?

How many & where: _____

Seal Date: _____

Y

N

Seal Name: _____

5. Did the custody seals arrive unbroken and intact upon arrival?

Y

N

6. COC#:

7. Were Custody papers filled out properly (ink,signed, legible, project information etc)?

Y

N

8. Were custody papers sealed in a plastic bag?

Y

N

9. Did you sign the COC in the appropriate place?

Y

N

10. Was enough ice used to chill the cooler?

Y N

Temp. of cooler: _____

0.3°C, 0.7°C

B. Log-In: Date samples were logged in:

EZ

By: 02/20/13

11. Were all bottles sealed in separate plastic bags?

Y

N

12. Did all bottles arrive unbroken and were labels in good condition?

Y

N

13. Were all bottle labels complete(ID,Date,time,etc.)

Y

N

14. Did all bottle labels agree with custody papers?

Y

N

15. Were the correct containers used for the tests indicated?

Y

N

16. Were samples received at the correct pH?

Y

N

17. Was sufficient amount of sample sent for the tests indicated?

Y

N

18. Were all samples submitted within holding time?

Y

N

19. Were all containers used within AEL's expiration date?**

Y

N

20. Were VOA samples absent of greater than pea-sized bubbles?

Y

N

(Note:Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

*If NO, List Sample ID's, Lab #s: _____

When bubbles are present in VOA samples they are labelled from smallest (or no bubbles) to largest. Lab to analyze VOA samples with no bubbles or
smallest bubbles first

20. Laboratory labeling verified by (initials): VB

Date: 2/20/13

**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.

February 27, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74874
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

- Case Narrative/Non-Conformance Summary
- Sample Log Sheet - Cover Page
- PCB Form 1 Data Sheet for Samples
- PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
- Chromatograms
- Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:

Sample SE-SB-219 (74874-10) was listed on the chain of custody (COC) but was not present in the cooler. The client was contacted and informed the laboratory that the sample was not collected.

PCBs by EPA Method 8082:

No results were reported below the quantitation limit.

Samples 74874-6 thru 74874-20 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument.

Sample 74873-18 had high recoveries for surrogate Tetrachloro-m-xylene. Decachlorobiphenyl surrogate recoveries were in control. Results were reported with a comment to this affect. Some samples had interference that prevented the determination of a surrogate recovery. Results were reported off the column without interferences when possible, or the sample was reported with a comment to this affect.

The MS/MSD analyzed on sample 74874-2 had low recoveries for PCB 1260. The parent sample had PCB 1260 detected. The laboratory control samples (L022113PSOX,RR/LD022113PSOX, RR) were in control. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC

Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
Summit Environmental Consultants Inc.
640 Main Street
Lewiston ME 04240

Report Number: 74874

Revision: Rev. 0

Re: MAINE ENERGY (Project No: 12-3259.1)

Enclosed are the results of the analyses on your sample(s). Samples were received on 20 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Sample Analysis: The attached pages detail the Client Sample IDs, Lab Sample IDs, and Analyses requested

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

If you have any questions on these results, please do not hesitate to contact us.

Authorized signature _____
Stephen L. Knollmeyer Lab. Director
Date _____

**This report shall not be reproduced, except in full, without the written
consent of Analytics Environmental Laboratory, LLC.**

CLIENT: Summit Environmental
 Consultants Inc.

REPORT NUMBER: 74874

REV: Rev. 0

PROJECT: MAINE ENERGY (Project No: 12-3259.1)

<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74874-1	02/19/13	SE-SB-221 (0-2')	EPA 8082 (PCBs only)	
74874-2	02/19/13	SE-SB-221 (2-4')	EPA 8082 (PCBs only)	
74874-3	02/19/13	SE-SB-221 (4-5')	EPA 8082 (PCBs only)	
74874-4	02/19/13	SE-SB-220 (0-2')	EPA 8082 (PCBs only)	
74874-5	02/19/13	SE-SB-220 (2-4')	EPA 8082 (PCBs only)	
74874-6	02/19/13	SE-SB-220 (4-6')	EPA 8082 (PCBs only)	
74874-7	02/19/13	SE-SB-220 (6-8')	EPA 8082 (PCBs only)	
74874-8	02/19/13	SE-SB-220 (8-10')	EPA 8082 (PCBs only)	
74874-9	02/19/13	SE-SB-220 (10-11')	EPA 8082 (PCBs only)	
74874-11	02/19/13	SE-SB-219 (0-2')	EPA 8082 (PCBs only)	
74874-12	02/19/13	SE-SB-219 (2-4')	EPA 8082 (PCBs only)	
74874-13	02/19/13	SE-SB-219 (4-6')	EPA 8082 (PCBs only)	
74874-14	02/19/13	SE-SB-219 (6-8')	EPA 8082 (PCBs only)	
74874-15	02/19/13	SE-SB-219 (8-10')	EPA 8082 (PCBs only)	
74874-16	02/19/13	SE-SB-224 (0-2')	EPA 8082 (PCBs only)	
74874-17	02/19/13	SE-SB-224 (2-4')	EPA 8082 (PCBs only)	
74874-18	02/19/13	SE-SB-224 (6-8')	EPA 8082 (PCBs only)	
74874-19	02/19/13	SE-SB-225 (0-2')	EPA 8082 (PCBs only)	
74874-20	02/19/13	SE-SB-225 (2-4')	EPA 8082 (PCBs only)	

Surrogate Compound Limits

	Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water				
1,4-Difluorobenzene		70-130		EPA 524.2
Bromofluorobenzene		70-130		
1,2-Dichlorobenzene-d4		70-130		
Volatile Organic Compounds				
1,2-Dichloroethane-d4		70-120	70-120	EPA 624/8260B
Toluene-d8		85-120	85-120	
Bromofluorobenzene		75-120	75-120	
Semi-Volatile Organic Compounds				
2-Fluorophenol		20-110	35-105	EPA 625/8270C
d5-Phenol		15-110	40-100	
d5-nitrobenzene		40-110	35-100	
2-Fluorobiphenyl		50-110	45-105	
2,4,6-Tribromophenol		40-110	40-125	
d14-p-terphenyl		50-130	30-125	
PAH's by SIM				
d5-nitrobenzene		21-110	35-110	EPA 8270C
2-Fluorobiphenyl		36-121	45-105	
d14-p-terphenyl		33-141	30-125	
Pesticides and PCBs				
2,4,5,6-Tetrachloro-m-xylene (TCX)		46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)		40-135	40-130	
Herbicides				
Dichloroacetic acid (DCAA)		30-150	30-150	
Gasoline Range Organics/TPH Gasoline				
Trifluorotoluene TFT (FID)		60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)		60-140	60-140	
Trifluorotoluene TFT (PID)		60-140	60-140	
Bromofluorobenzene (BFB) (PID)		60-140	60-140	
Diesel Range Organics/TPH Diesel				
m-terphenyl		60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons				
2,5-Dibromotoluene (PID)		70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)		70-130	70-130	
Extricable Petroleum Hydrocarbons				
1-chloro-octadecane (aliphatic)		40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)		40-140	40-140	
2-Fluorobiphenyl (Fractionation)		40-140	40-140	
2-Bromonaphthalene (fractionation)		40-140	40-140	

**PCB
DATA SUMMARIES**

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-221 (0-2')

Lab Sample ID: 74874-1
Matrix: Solid
Percent Solid: 73
Dilution Factor: 1.4
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	46	U
PCB-1221	46	U
PCB-1232	46	U
PCB-1242	46	U
PCB-1248	46	U
PCB-1254	46	U
PCB-1260	46	47

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	81	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG:

GC Column #1: STX-CLPesticides I

Sample: 74874-1,,A/C

Column ID: 0.25 mm

Data File: M67433.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	47	34	32.2	

Column to be used to flag RPD values greater than QC limit of 40%

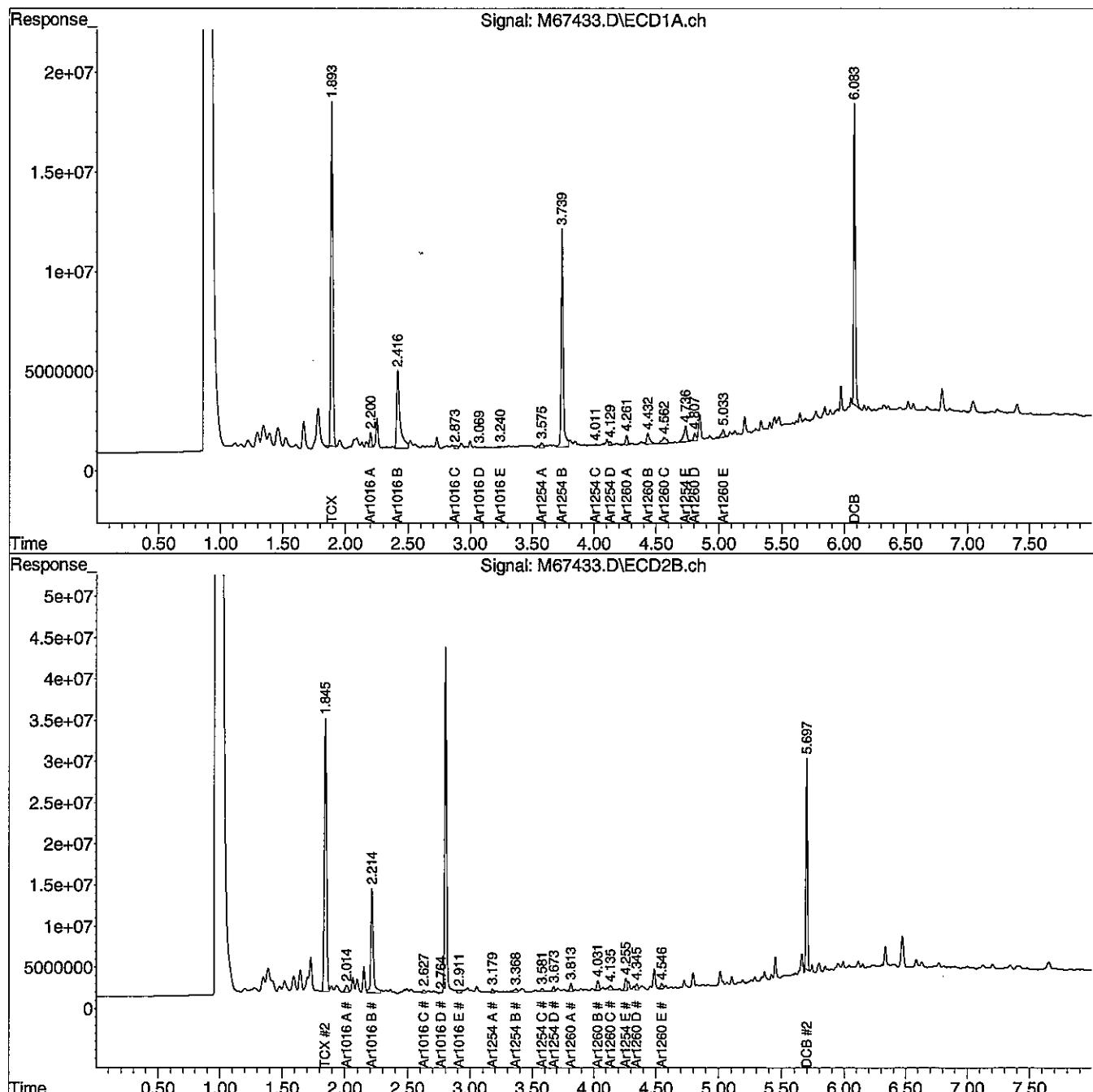
* Values outside QC limits

Comments: _____

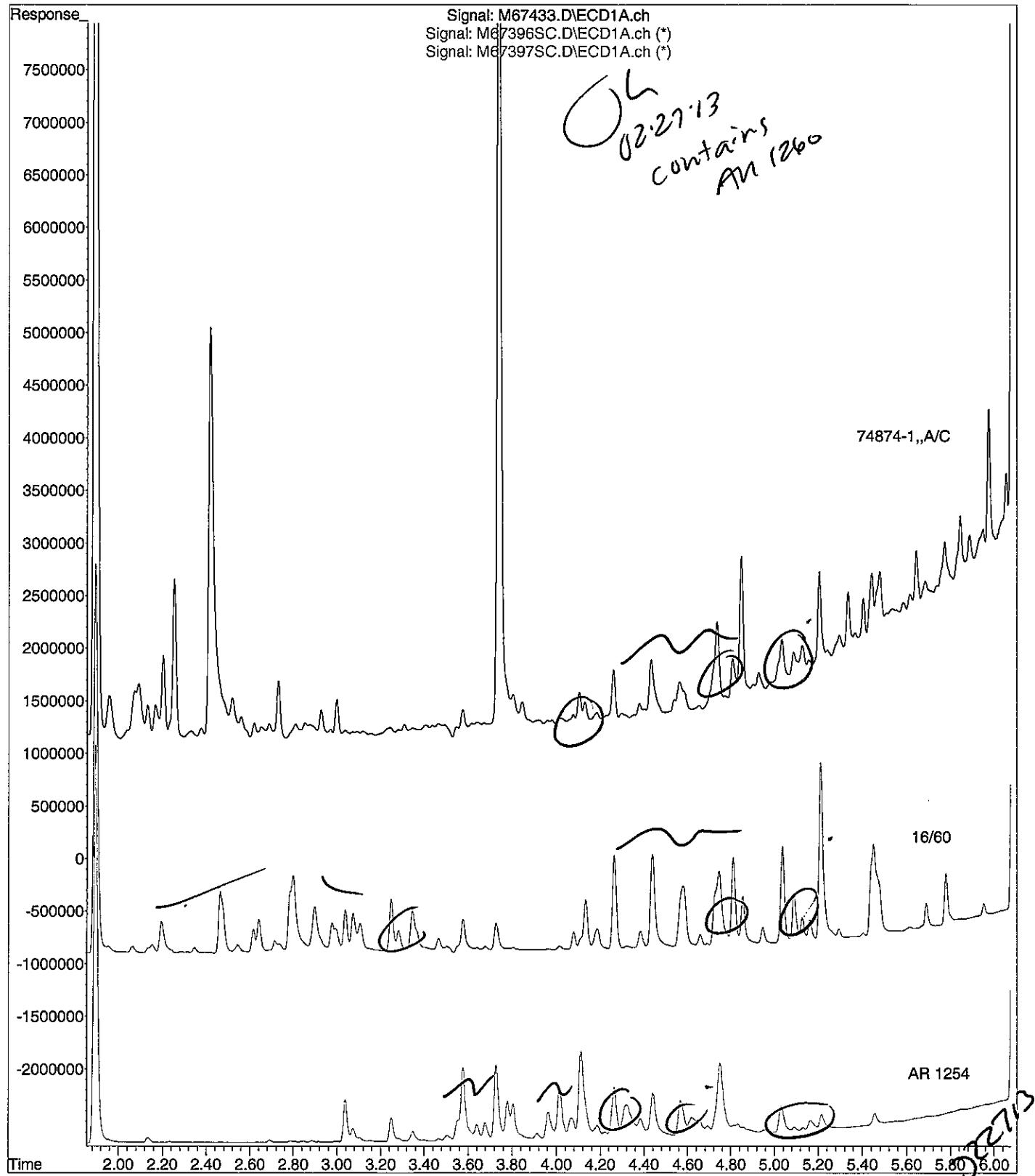
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67433.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 10:14 pm
 Operator : JK
 Sample : 74874-1,,A/C
 Misc : SOIL
 ALS Vial : 21 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:03:50 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67433.D
Operator : JK
Acquired : 26 Feb 2013 10:14 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74874-1,,A/C
Misc Info : SOIL
Vial Number: 21



Mr. John Cressey
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 Lewiston ME 04240

February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-221 (2-4')

Lab Sample ID:	74874-2
Matrix:	Solid
Percent Solid:	88
Dilution Factor:	1.1
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	36	U
PCB-1221	36	U
PCB-1232	36	U
PCB-1242	36	U
PCB-1248	36	U
PCB-1254	36	U
PCB-1260	36	308

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	102	%
Decachlorobiphenyl	71	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-2,,A/C

Column ID: 0.25 mm

Data File: M67434.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	308	268	13.9	

Column to be used to flag RPD values greater than QC limit of 40%

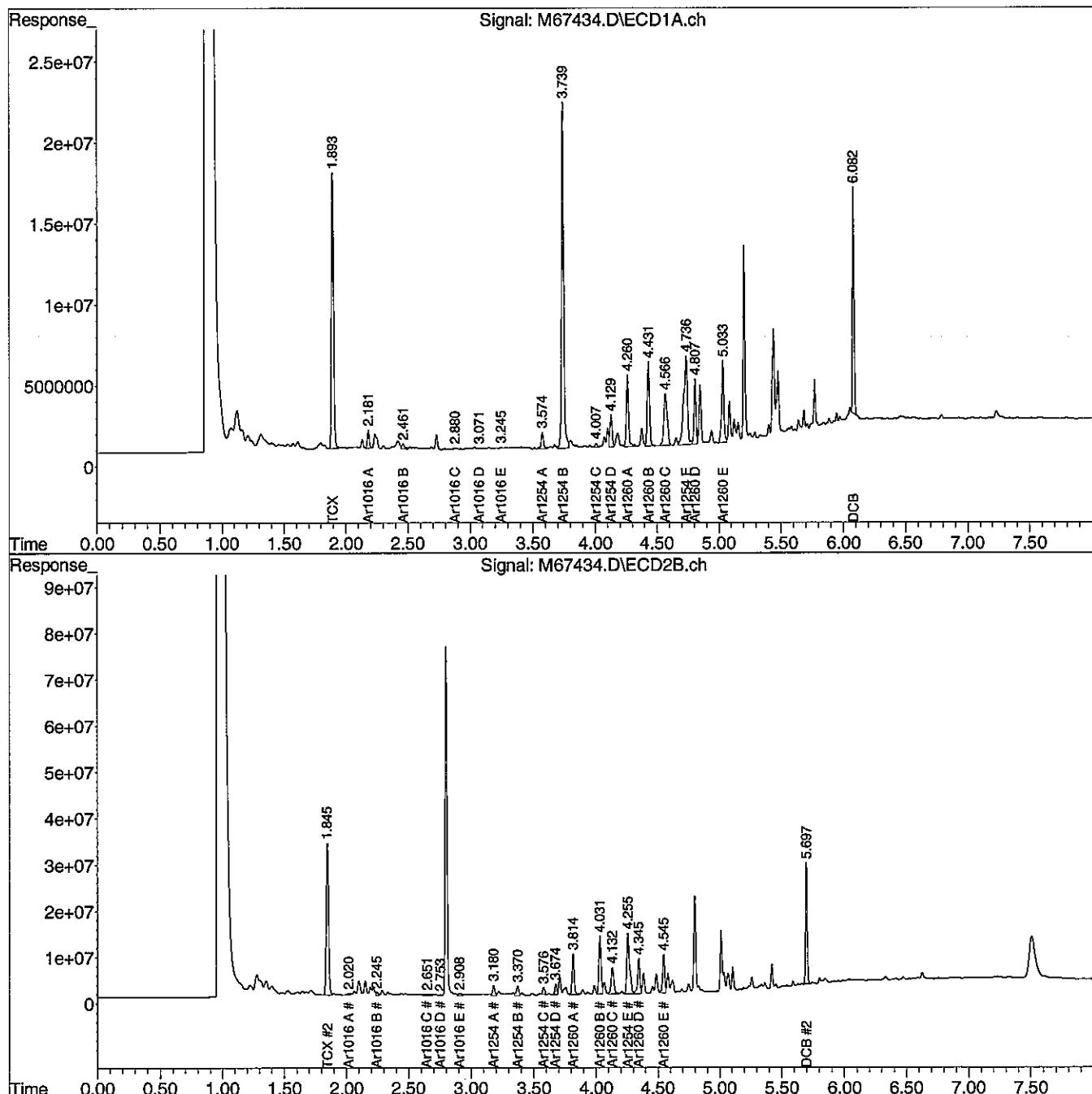
* Values outside QC limits

Comments: _____

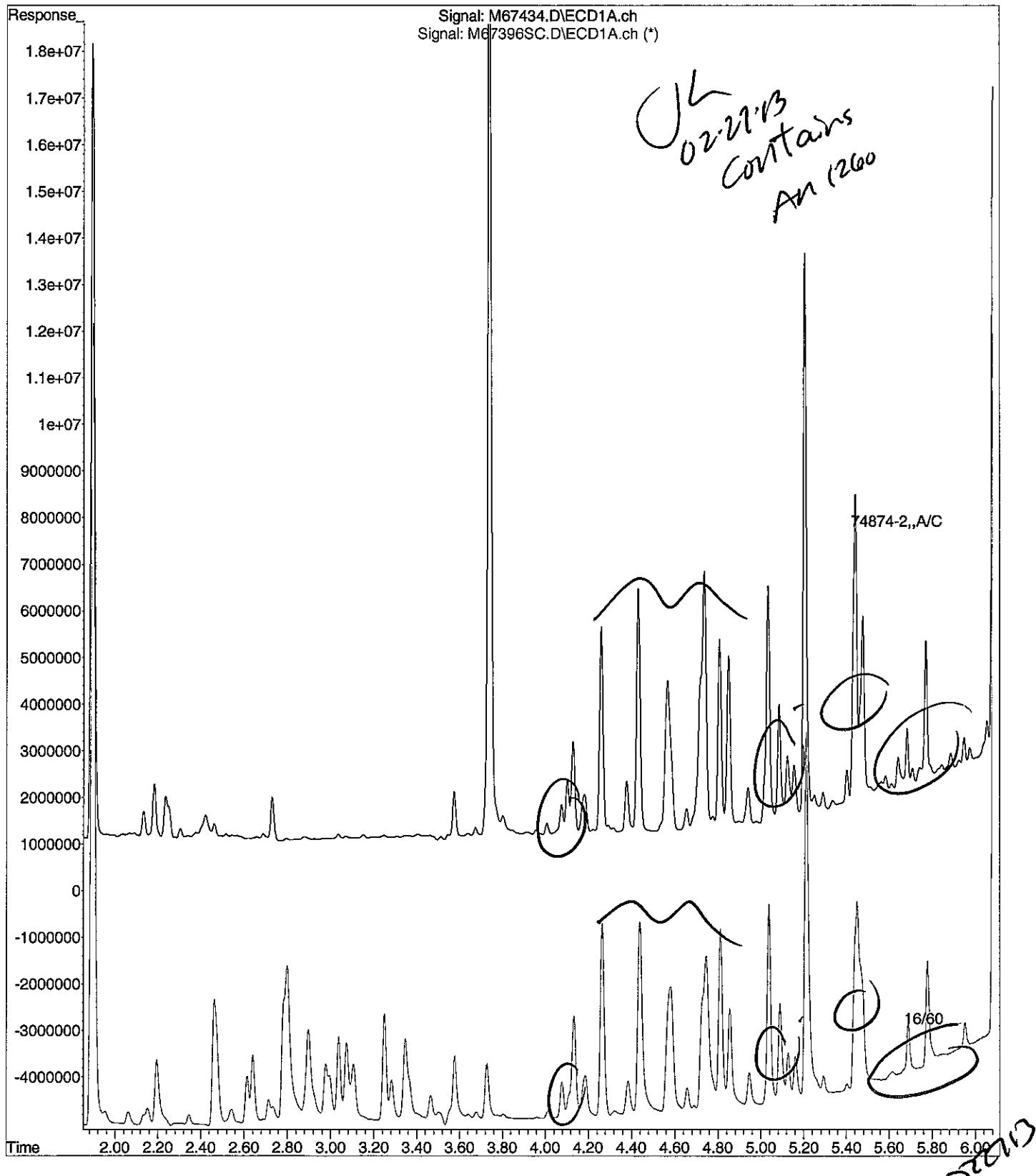
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67434.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 10:24 pm
 Operator : JK
 Sample : 74874-2,,A/C
 Misc : SOIL
 ALS Vial : 22 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:30 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67434.D
Operator : JK
Acquired : 26 Feb 2013 10:24 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74874-2,,A/C
Misc Info : SOIL
Vial Number: 22



Mr. John Cressey
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 640 Main Street
 Lewiston ME 04240

February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-221 (4-5')

Lab Sample ID: 74874-3
Matrix: Solid
Percent Solid: 85
Dilution Factor: 1.2
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	40	U
PCB-1221	40	U
PCB-1232	40	U
PCB-1242	40	U
PCB-1248	40	U
PCB-1254	40	U
PCB-1260	40	68

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	106	%
Decachlorobiphenyl	71	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-3,,A/C

Column ID: 0.25 mm

Data File: M67437.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.2

Column ID: 0.25 mm

	Column #1	Column #2	RPD	#
COMPOUND	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	64	68	6.7	

Column to be used to flag RPD values greater than QC limit of 40%

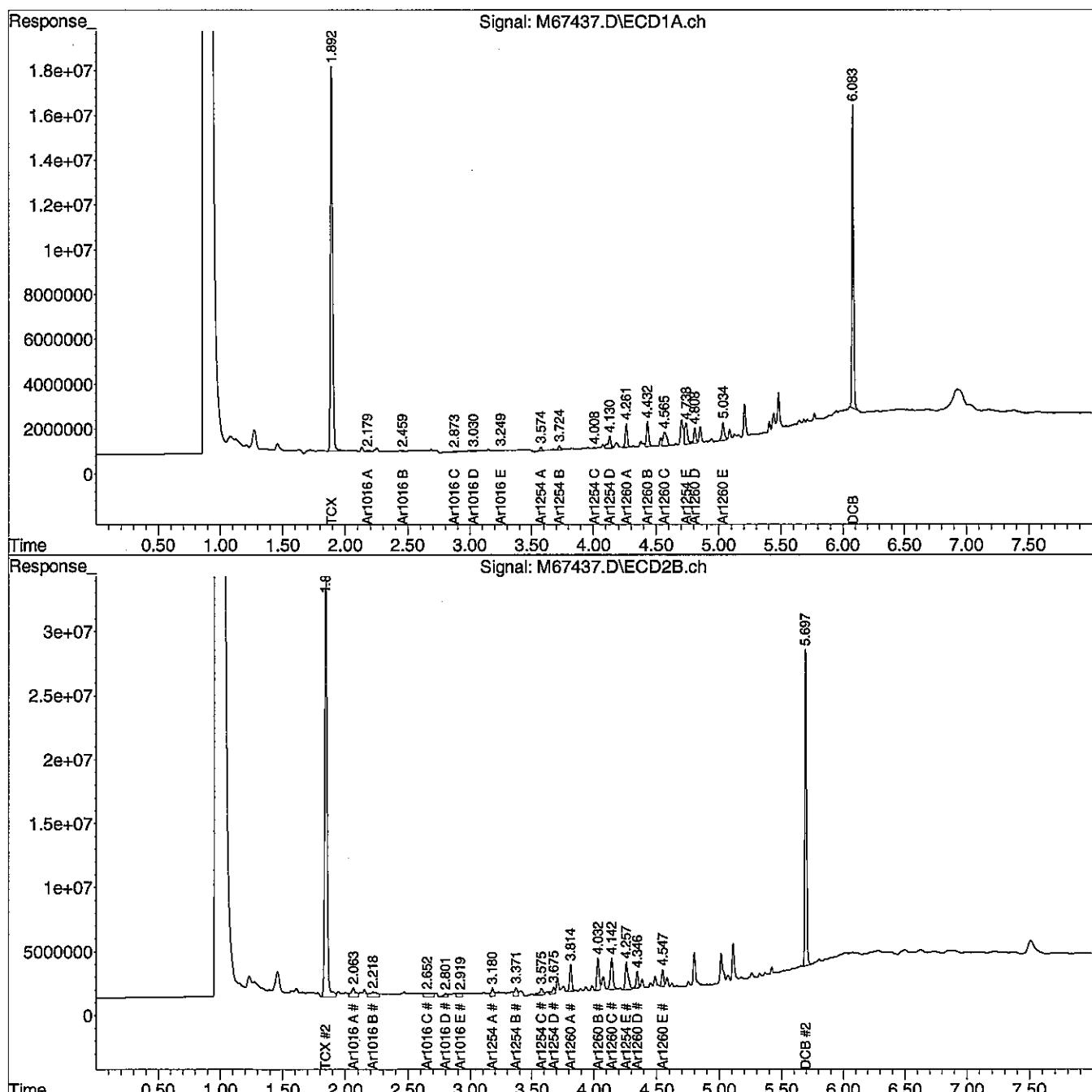
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67437.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 10:54 pm
 Operator : JK
 Sample : 74874-3,,A/C
 Misc : SOIL
 ALS Vial : 25 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:09:17 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-220 (0-2')

Lab Sample ID:	74874-4
Matrix:	Solid
Percent Solid:	73
Dilution Factor:	1.3
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	43	U
PCB-1221	43	U
PCB-1232	43	U
PCB-1242	43	U
PCB-1248	43	U
PCB-1254	43	U
PCB-1260	43	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	105	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

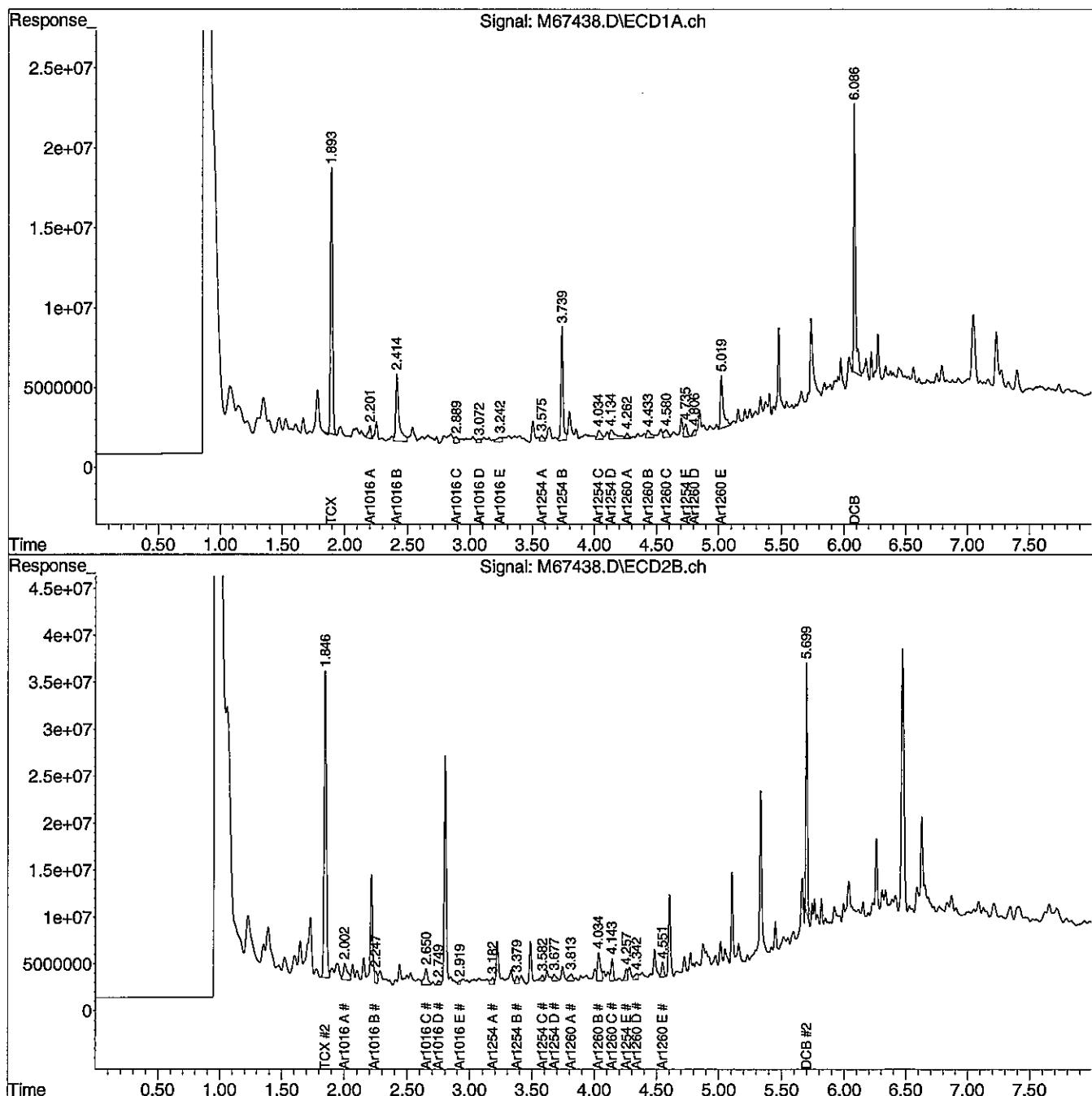
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67438.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:04 pm
 Operator : JK
 Sample : 74874-4,,A/C
 Misc : SOIL
 ALS Vial : 26 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:38 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-220 (2-4')

Lab Sample ID: 74874-5
Matrix: Solid
Percent Solid: 88
Dilution Factor: 1.0
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	284

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	99	%
Decachlorobiphenyl	78	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-5,,A/C

Column ID: 0.25 mm

Data File: M67439.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD #
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)	
PCB 1260	284	248	13.5

Column to be used to flag RPD values greater than QC limit of 40%

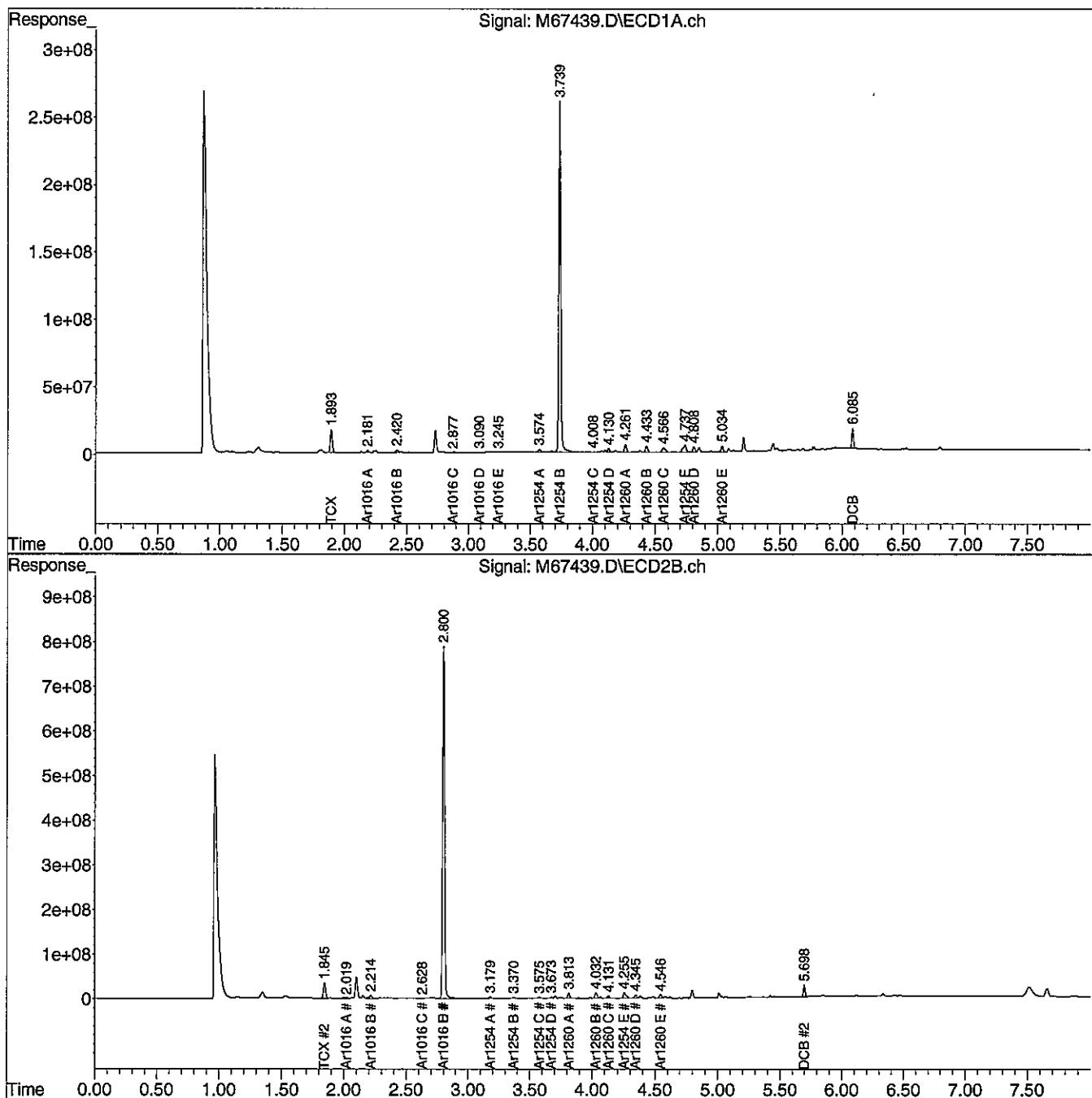
* Values outside QC limits

Comments: _____

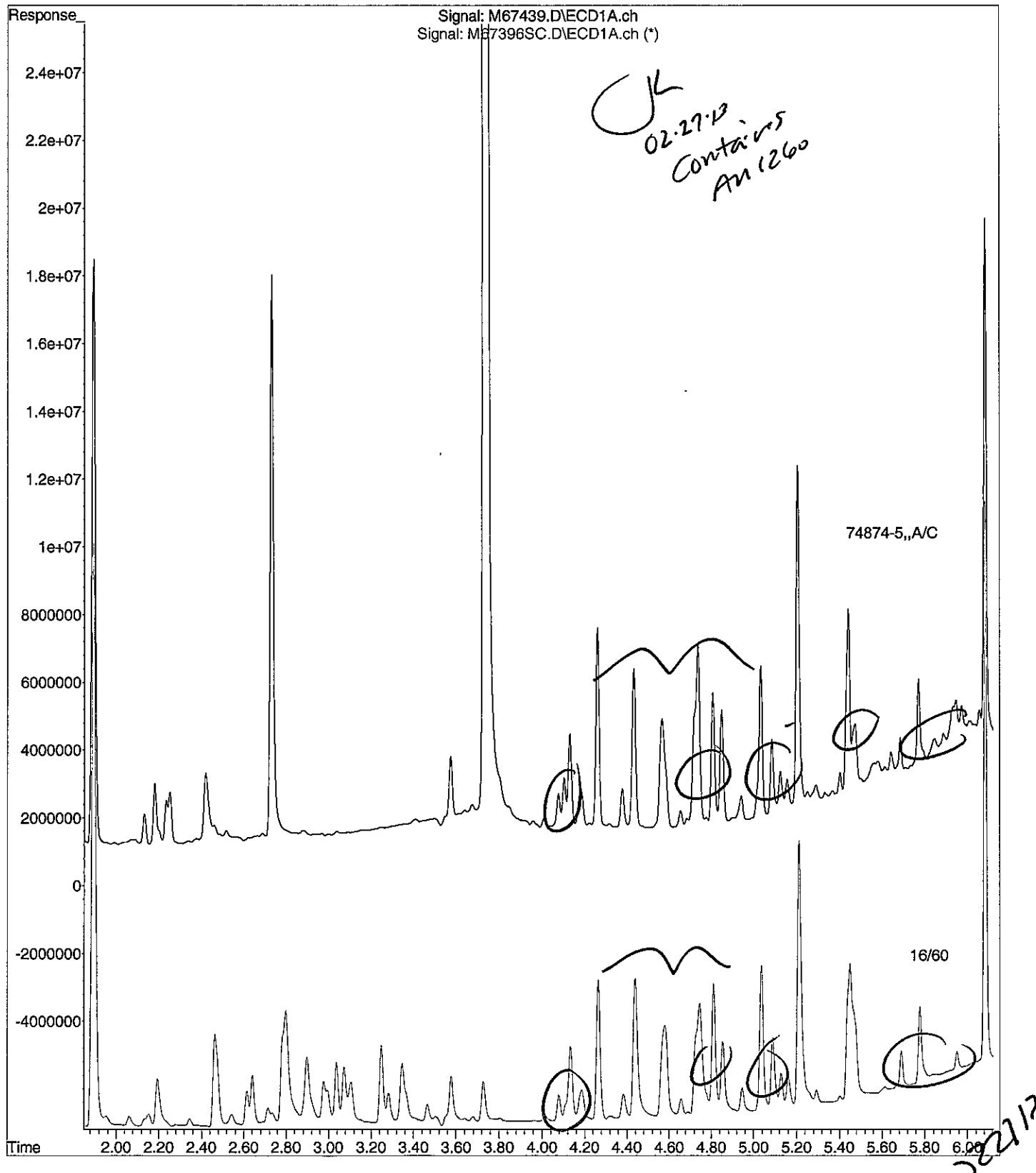
Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67439.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:15 pm
 Operator : JK
 Sample : 74874-5, ,A/C
 Misc : SOIL
 ALS Vial : 27 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:40 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022613-M\M67439.D
Operator : JK
Acquired : 26 Feb 2013 11:15 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74874-5,,A/C
Misc Info : SOIL
Vial Number: 27



Mr. John Cressey
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 Lewiston ME 04240

February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-220 (4-6')

Lab Sample ID: 74874-6
Matrix: Solid
Percent Solid: 90
Dilution Factor: 5
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	2610

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	87	%	
Decachlorobiphenyl	71	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-6,1:5,A/C

Column ID: 0.25 mm

Data File: M67440.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2611	2183	17.9	

Column to be used to flag RPD values greater than QC limit of 40%

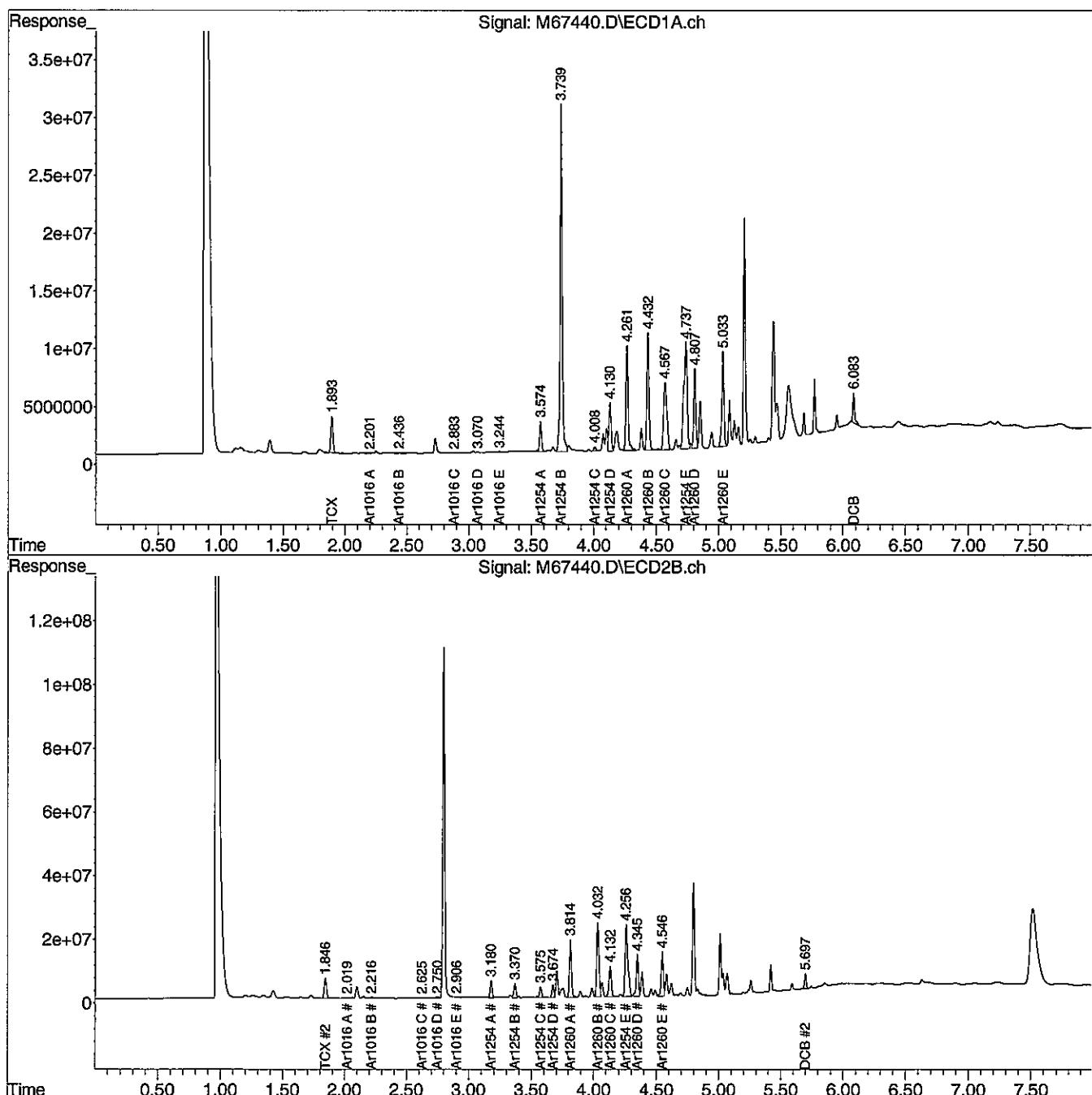
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67440.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:25 pm
 Operator : JK
 Sample : 74874-6,1:5,,A/C
 Misc : SOIL
 ALS Vial : 28 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:42 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-220 (6-8")

Lab Sample ID: 74874-7
Matrix: Solid
Percent Solid: 92
Dilution Factor: 22
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	726	U
PCB-1221	726	U
PCB-1232	726	U
PCB-1242	726	U
PCB-1248	726	U
PCB-1254	726	U
PCB-1260	726	13900

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-7,1:20,,A/C

Column ID: 0.25 mm

Data File: M67441.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 21.7

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD #
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)	
PCB 1260	13895	11507	18.8

Column to be used to flag RPD values greater than QC limit of 40%

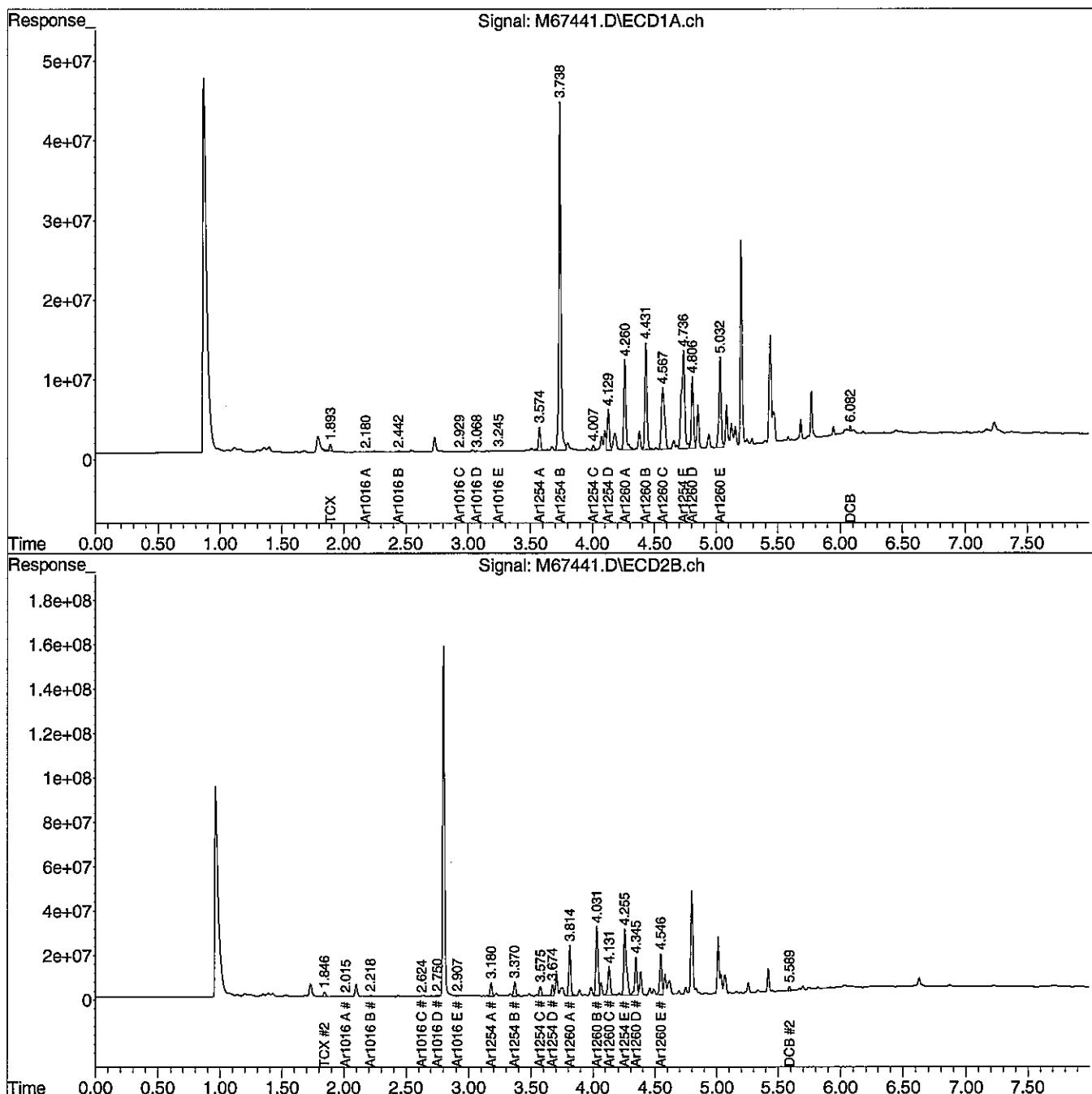
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67441.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:35 pm
 Operator : JK
 Sample : 74874-7,1:20,,A/C
 Misc : SOIL
 ALS Vial : 29 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:44 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-220 (8-10')

Lab Sample ID: 74874-8
Matrix: Solid
Percent Solid: 86
Dilution Factor: 27
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	891	U
PCB-1221	891	U
PCB-1232	891	U
PCB-1242	891	U
PCB-1248	891	U
PCB-1254	891	U
PCB-1260	891	19800

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG:

GC Column #1: STX-CLPesticides I

Sample: 74874-8,1:25,,A/C

Column ID: 0.25 mm

Data File: M67442.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 27.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	19812	16914	15.8	

Column to be used to flag RPD values greater than QC limit of 40%

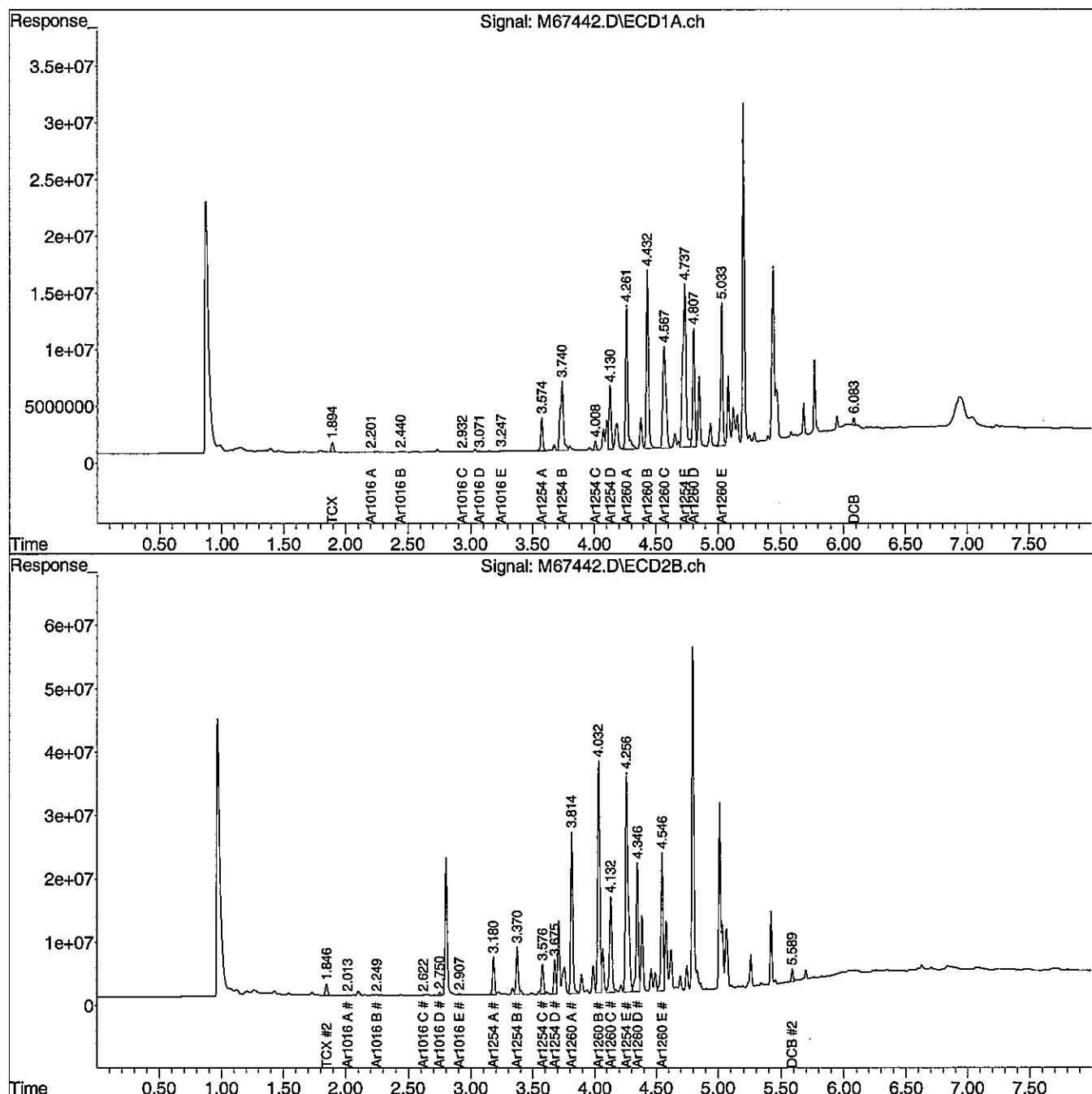
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67442.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:45 pm
 Operator : JK
 Sample : 74874-8,1:25,,A/C
 Misc : SOIL
 ALS Vial : 30 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:46 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY

Lab Sample ID: 74874-9

Project Number: 12-3259.1

Matrix: Solid

Field Sample ID: SE-SB-220 (10-11')

Percent Solid: 82

Dilution Factor: 609

Collection Date: 02/19/13

Lab Receipt Date: 02/20/13

Extraction Date: 02/21/13

Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	20100	U
PCB-1221	20100	U
PCB-1232	20100	U
PCB-1242	20100	U
PCB-1248	20100	U
PCB-1254	20100	U
PCB-1260	20100	489000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene * %

Decachlorobiphenyl * %

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-9,1:500,,A/C

Column ID: 0.25 mm

Data File: M67443.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 609.0

Column ID: 0.25 mm

	Column #1	Column #2	RPD	#
COMPOUND	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	489243	420718	15.1	

Column to be used to flag RPD values greater than QC limit of 40%

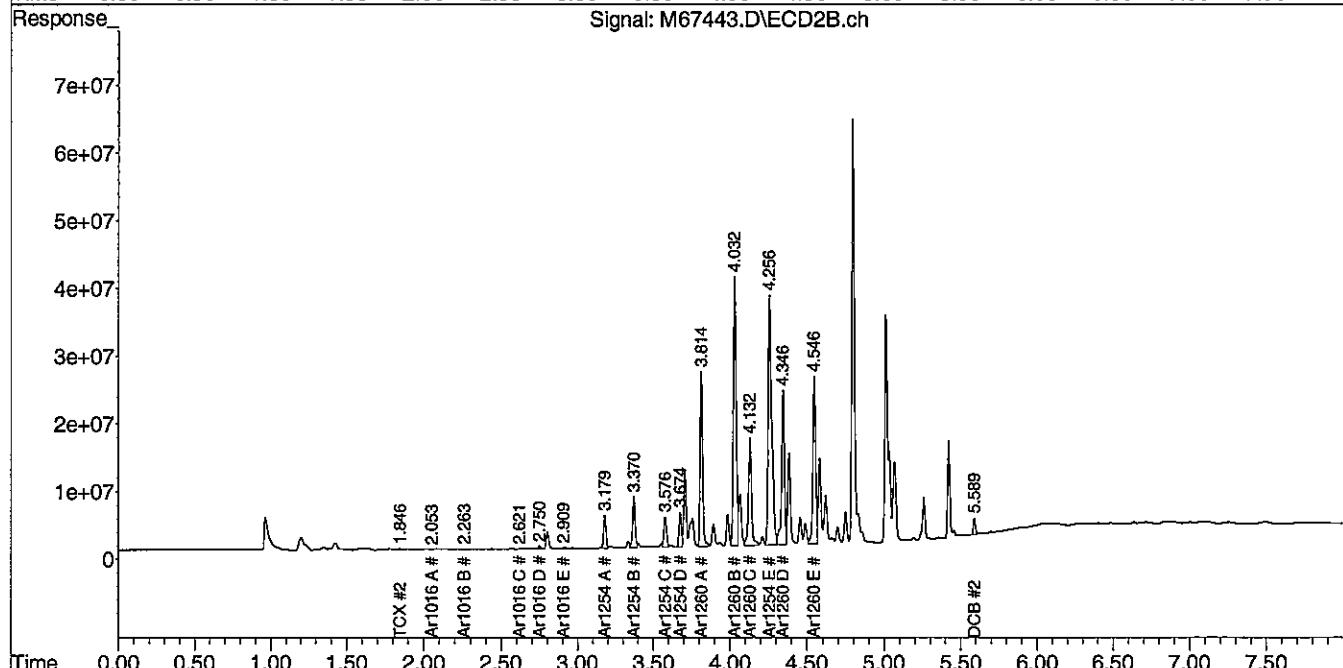
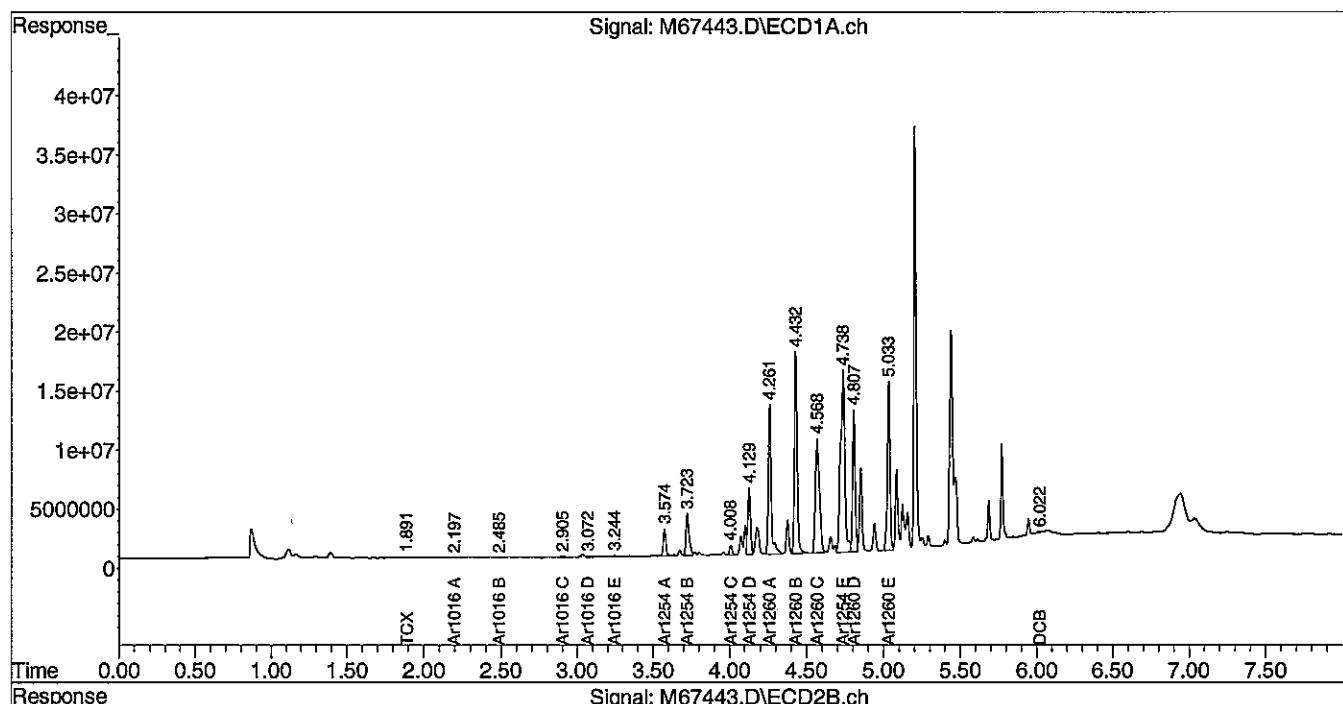
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67443.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 11:55 pm
 Operator : JK
 Sample : 74874-9,1:500,,A/C
 Misc : SOIL
 ALS Vial : 31 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:48 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 26, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-219 (0-2")

Lab Sample ID:	74874-11
Matrix:	Solid
Percent Solid:	92
Dilution Factor:	5
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/23/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	118	%
Decachlorobiphenyl	107	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

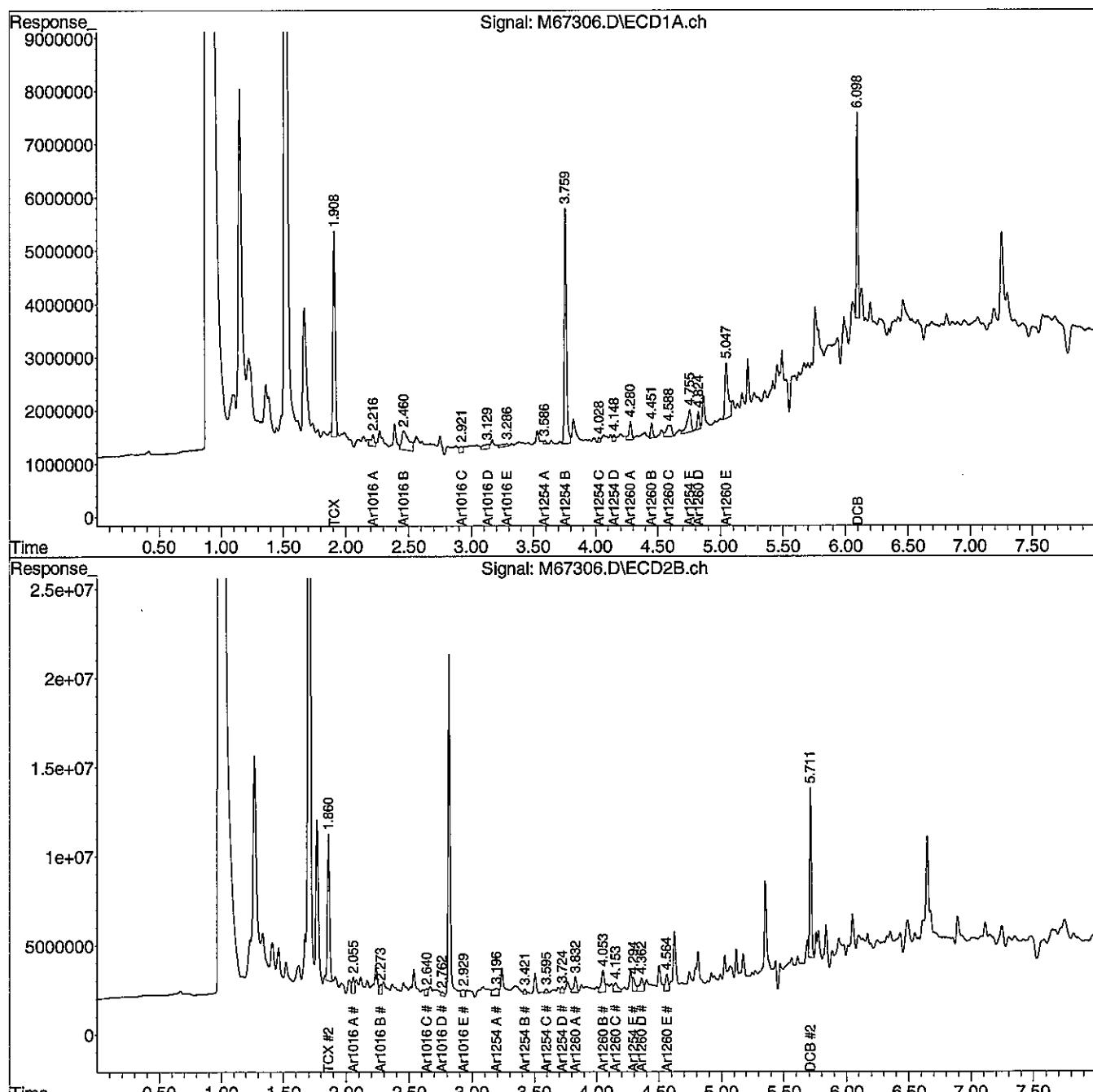
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * Quantitation limits increased due to the sample matrix affect.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67306.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 23 Feb 2013 6:36 am
 Operator : JK
 Sample : 74874-11,1:5,,A/C
 Misc : SOIL
 ALS Vial : 42 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 15:57:17 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 26, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-219 (2-4')

Lab Sample ID:	74874-12
Matrix:	Solid
Percent Solid:	92
Dilution Factor:	5
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/23/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	165	U
PCB-1221	165	U
PCB-1232	165	U
PCB-1242	165	U
PCB-1248	165	U
PCB-1254	165	U
PCB-1260	165	380

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	125	%
Decachlorobiphenyl	115	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-12,1:5,,A/C

Column ID: 0.25 mm

Data File: M67307.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.1

Column ID: 0.25 mm

COMPOUND	Column #1		Column #2		#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)	RPD		
PCB 1260	380	372	2.0		

Column to be used to flag RPD values greater than QC limit of 40%

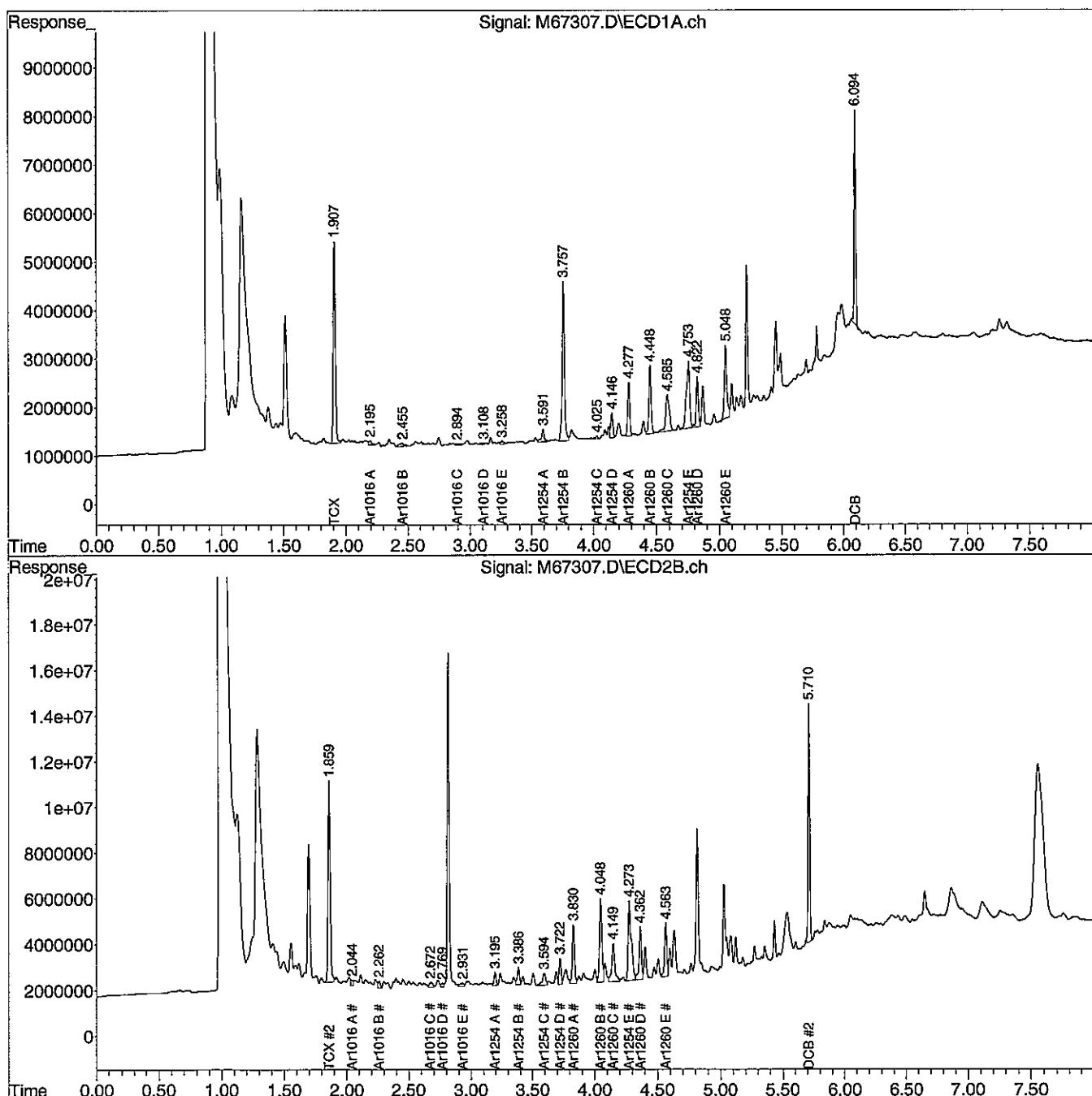
* Values outside QC limits

Comments: _____

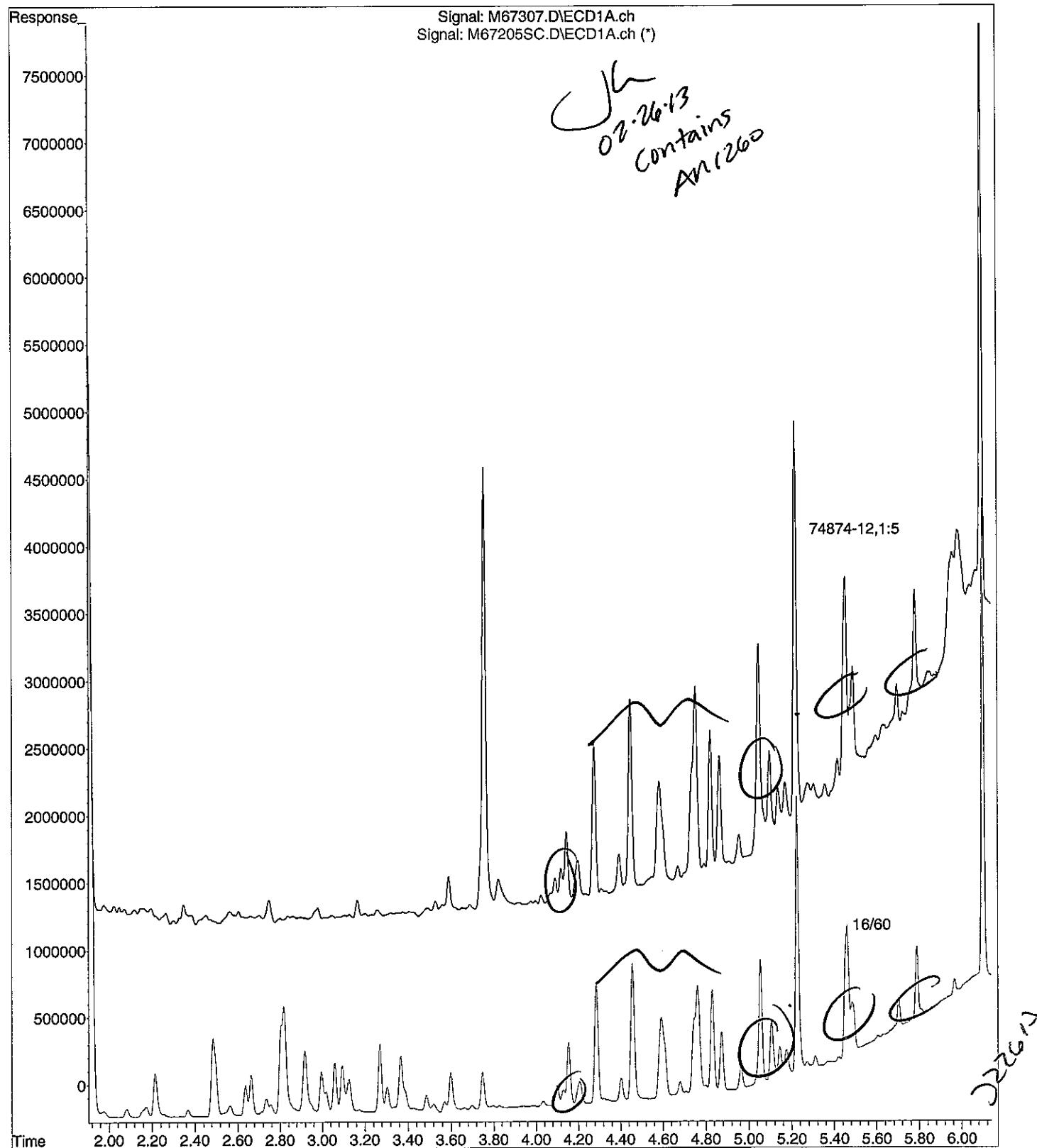
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67307.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 23 Feb 2013 6:47 am
 Operator : JK
 Sample : 74874-12,1:5,,A/C
 Misc : SOIL
 ALS Vial : 43 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 16:09:56 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67307.D
Operator : JK
Acquired : 23 Feb 2013 6:47 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74874-12,1:5,,A/C
Misc Info : SOIL
Vial Number: 43



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-219 (4-6')

Lab Sample ID:	74874-13
Matrix:	Solid
Percent Solid:	66
Dilution Factor:	15
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	495	U
PCB-1221	495	U
PCB-1232	495	U
PCB-1242	495	U
PCB-1248	495	U
PCB-1254	495	U
PCB-1260	495	11000

Surrogate Standard Recovery			
2,4,5,6-Tetrachloro-m-xylene	94	%	
Decachlorobiphenyl	95	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-13,1:10,,A/C

Column ID: 0.25 mm

Data File: M67444.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 15.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	10955	8644	23.6	

Column to be used to flag RPD values greater than QC limit of 40%

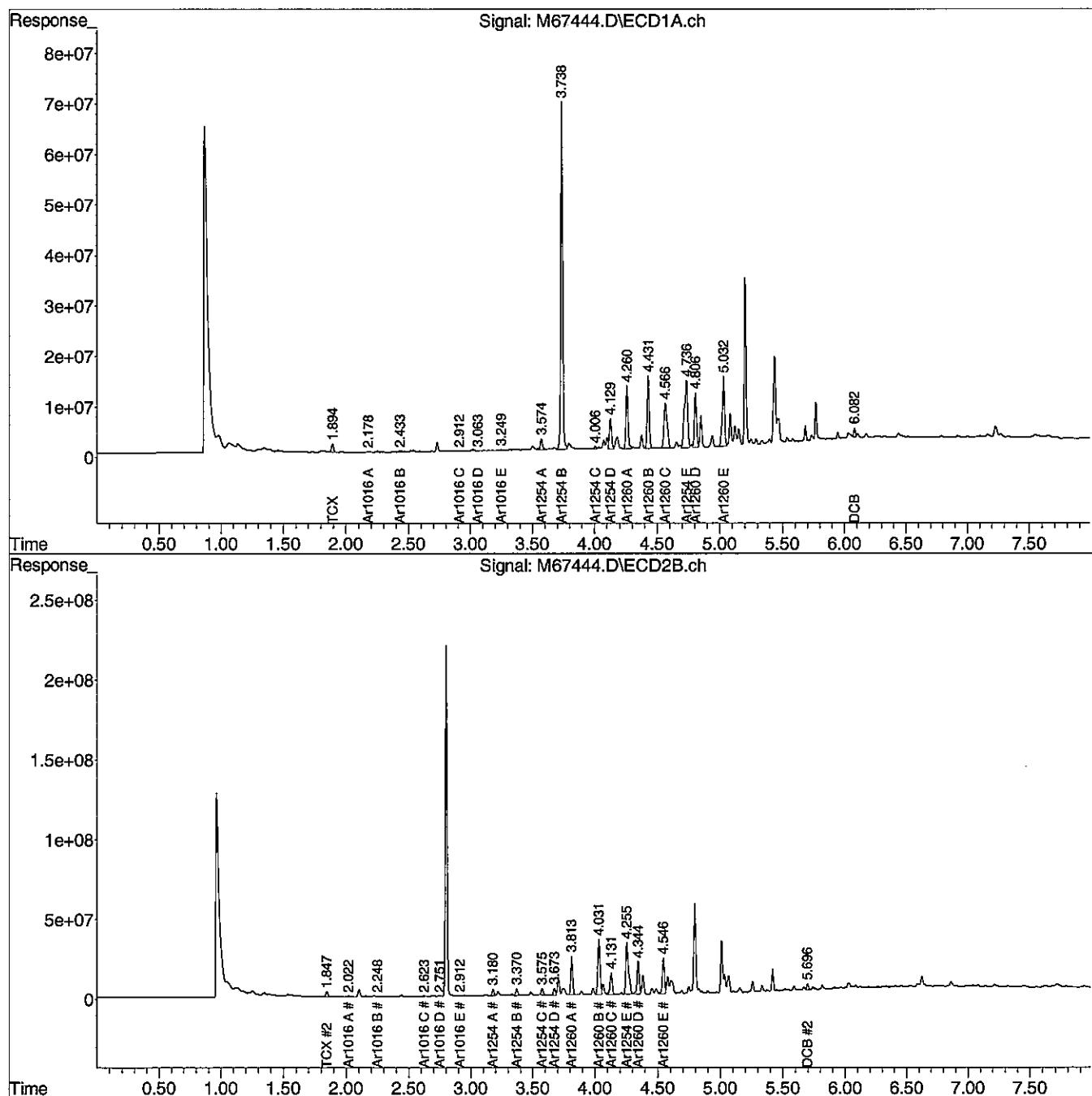
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67444.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:05 am
 Operator : JK
 Sample : 74874-13,1:10,,A/C
 Misc : SOIL
 ALS Vial : 32 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 12:15:11 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-219 (6-8")

Lab Sample ID:	74874-14
Matrix:	Solid
Percent Solid:	87
Dilution Factor:	54
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	1780	U
PCB-1221	1780	U
PCB-1232	1780	U
PCB-1242	1780	U
PCB-1248	1780	U
PCB-1254	1780	U
PCB-1260	1780	39800

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG:

GC Column #1: STX-CLPesticides I

Sample: 74874-14,1:50,,A/C

Column ID: 0.25 mm

Data File: M67445.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 53.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	39845	34549	14.2	

Column to be used to flag RPD values greater than QC limit of 40%

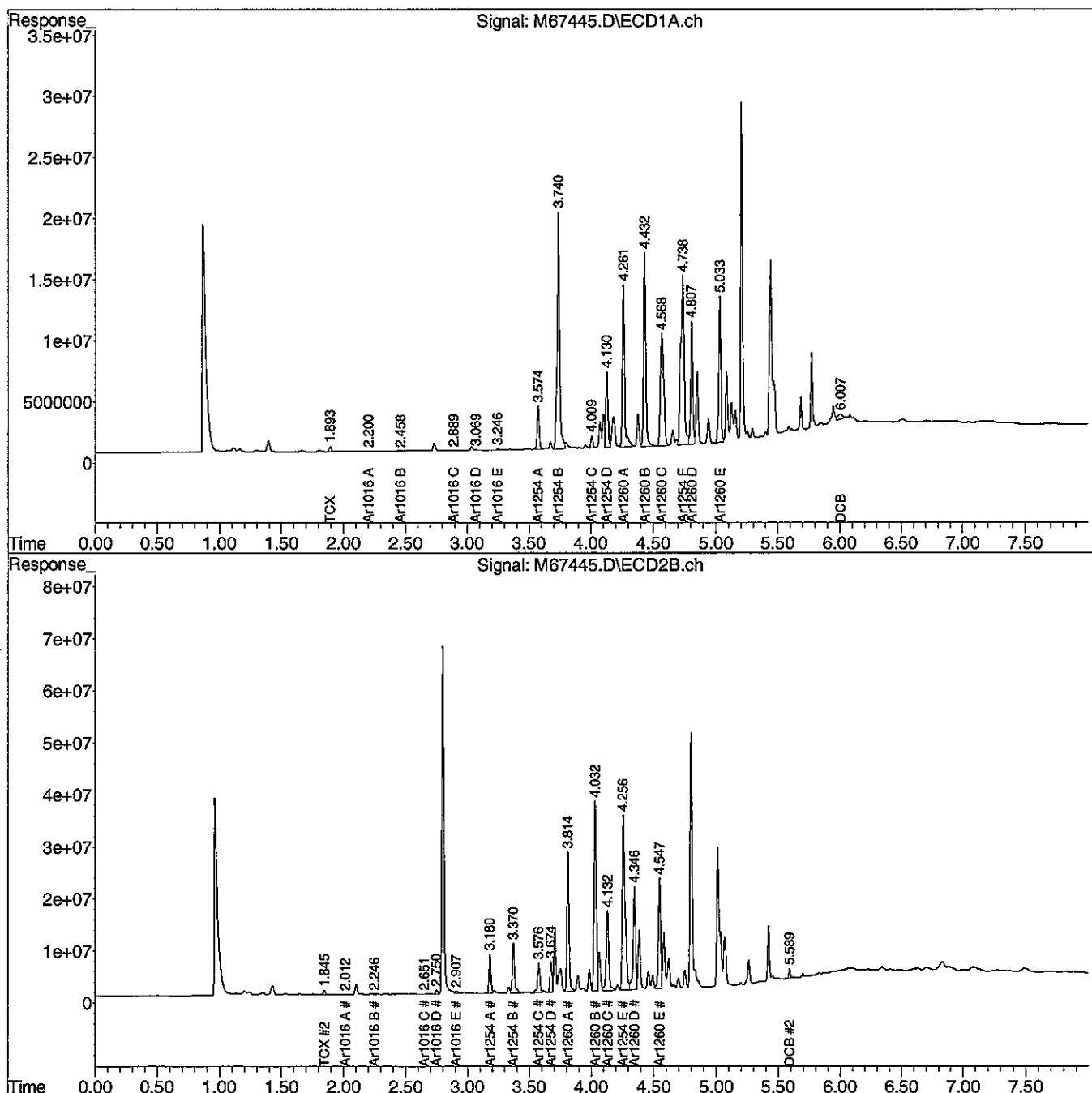
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67445.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:15 am
 Operator : JK
 Sample : 74874-14,1:50,,A/C
 Misc : SOIL
 ALS Vial : 33 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:52 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-219 (8-10')

Lab Sample ID: 74874-15
Matrix: Solid
Percent Solid: 91
Dilution Factor: 109
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	3600	U
PCB-1221	3600	U
PCB-1232	3600	U
PCB-1242	3600	U
PCB-1248	3600	U
PCB-1254	3600	U
PCB-1260	3600	54100

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-15,1:100,,A/C

Column ID: 0.25 mm

Data File: M67446.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 108.9

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	54112	47749	12.5	

Column to be used to flag RPD values greater than QC limit of 40%

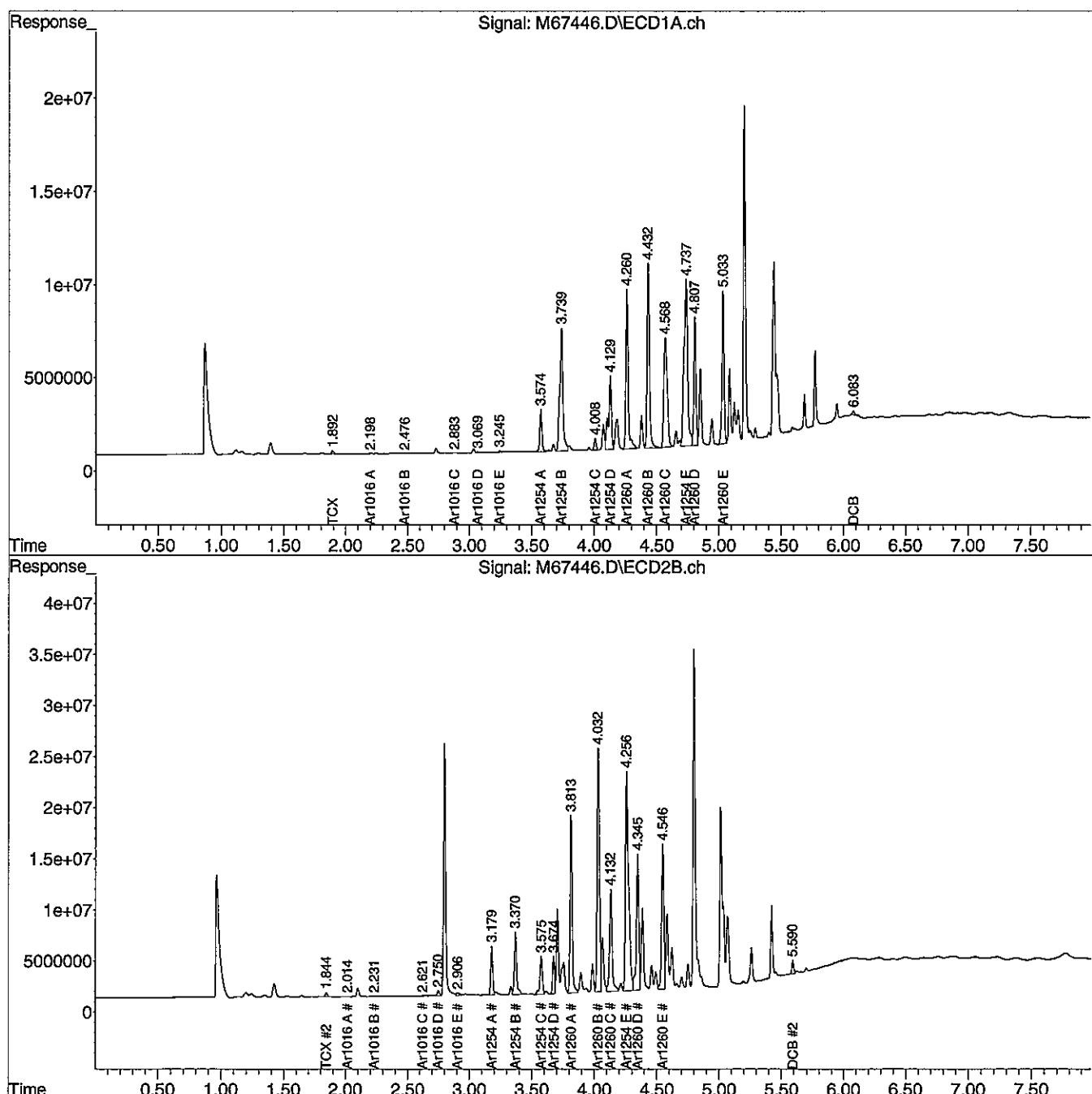
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67446.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:25 am
 Operator : JK
 Sample : 74874-15,1:100,,A/C
 Misc : SOIL
 ALS Vial : 34 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:54 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-224 (0-2")

Lab Sample ID: 74874-16
Matrix: Solid
Percent Solid: 89
Dilution Factor: 11
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	9550

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	117	%	
Decachlorobiphenyl	74	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-16,1:10,,A/C

Column ID: 0.25 mm

Data File: M67447.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.8

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	9550	8315	13.8	

Column to be used to flag RPD values greater than QC limit of 40%

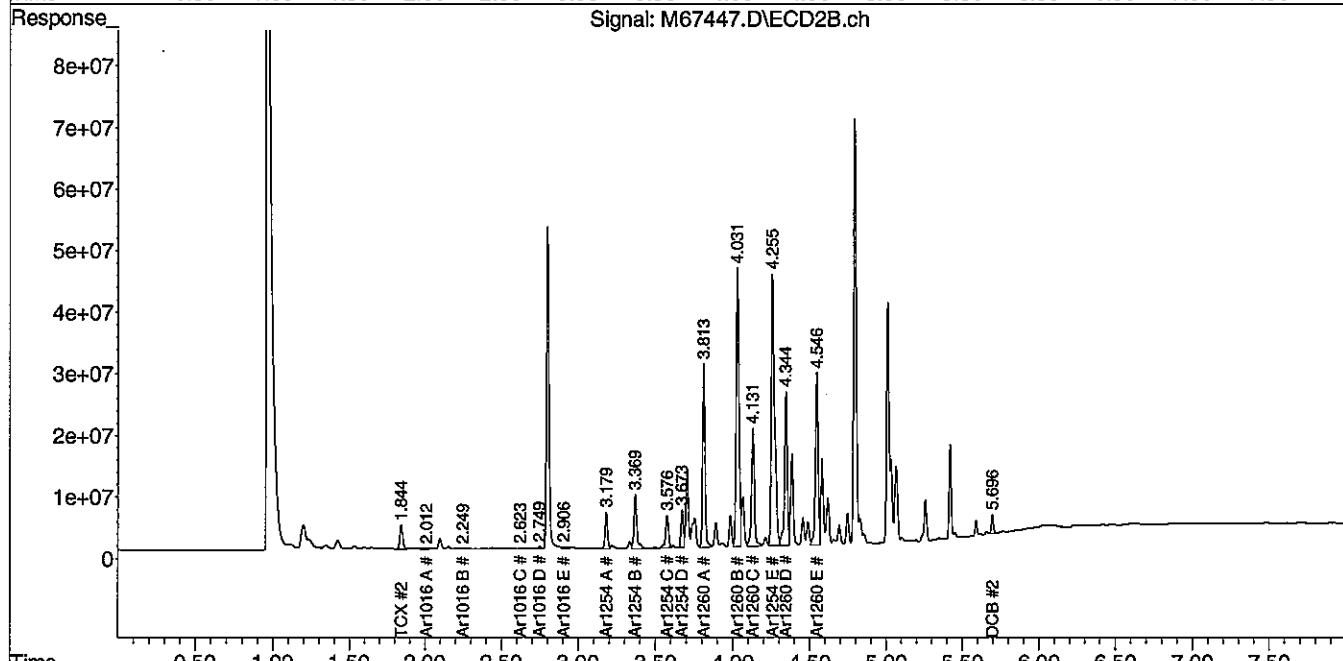
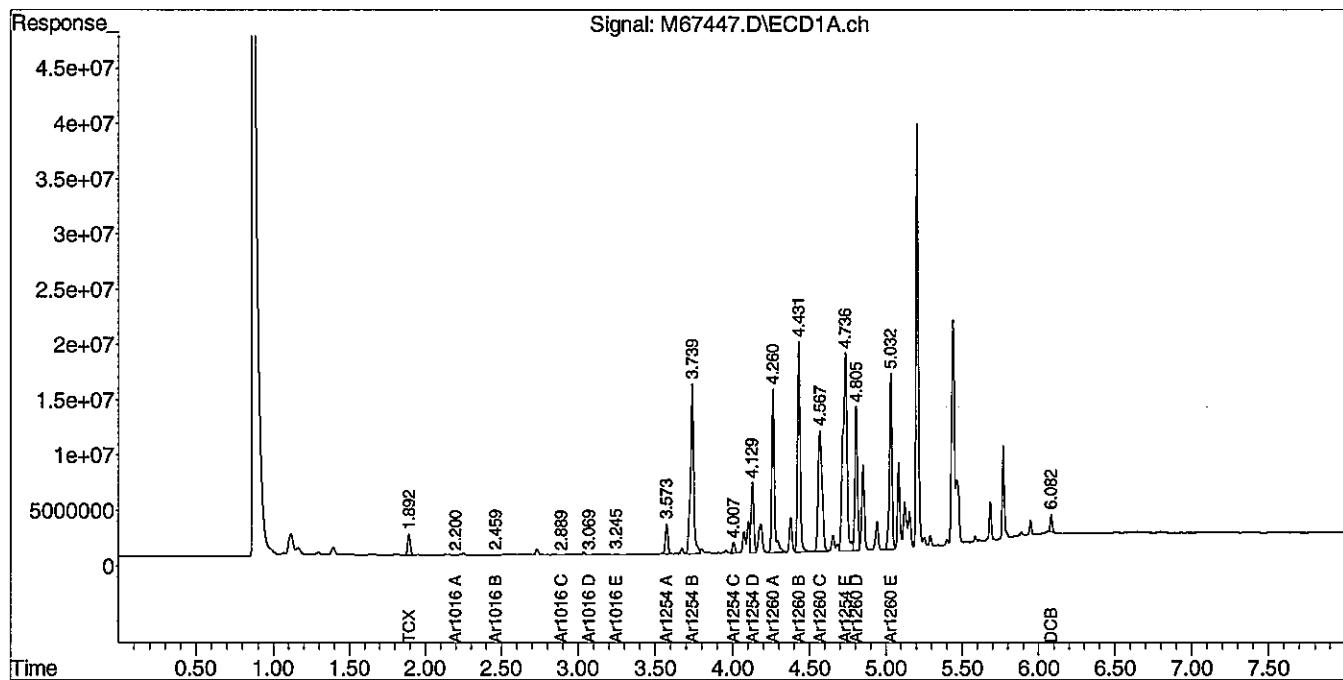
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67447.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:35 am
 Operator : JK
 Sample : 74874-16,1:10,,A/C
 Misc : SOIL
 ALS Vial : 35 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:56 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



Mr. John Cressey
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 640 Main Street
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February 26, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-224 (2-4")

Lab Sample ID: 74874-17
Matrix: Solid
Percent Solid: 84
Dilution Factor: 1.2
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/23/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	40	U
PCB-1221	40	U
PCB-1232	40	U
PCB-1242	40	U
PCB-1248	40	U
PCB-1254	40	U
PCB-1260	40	229

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	94	%
Decachlorobiphenyl	67	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-17,,A/C

Column ID: 0.25 mm

Data File: M67312.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	229	211	8.3	

Column to be used to flag RPD values greater than QC limit of 40%

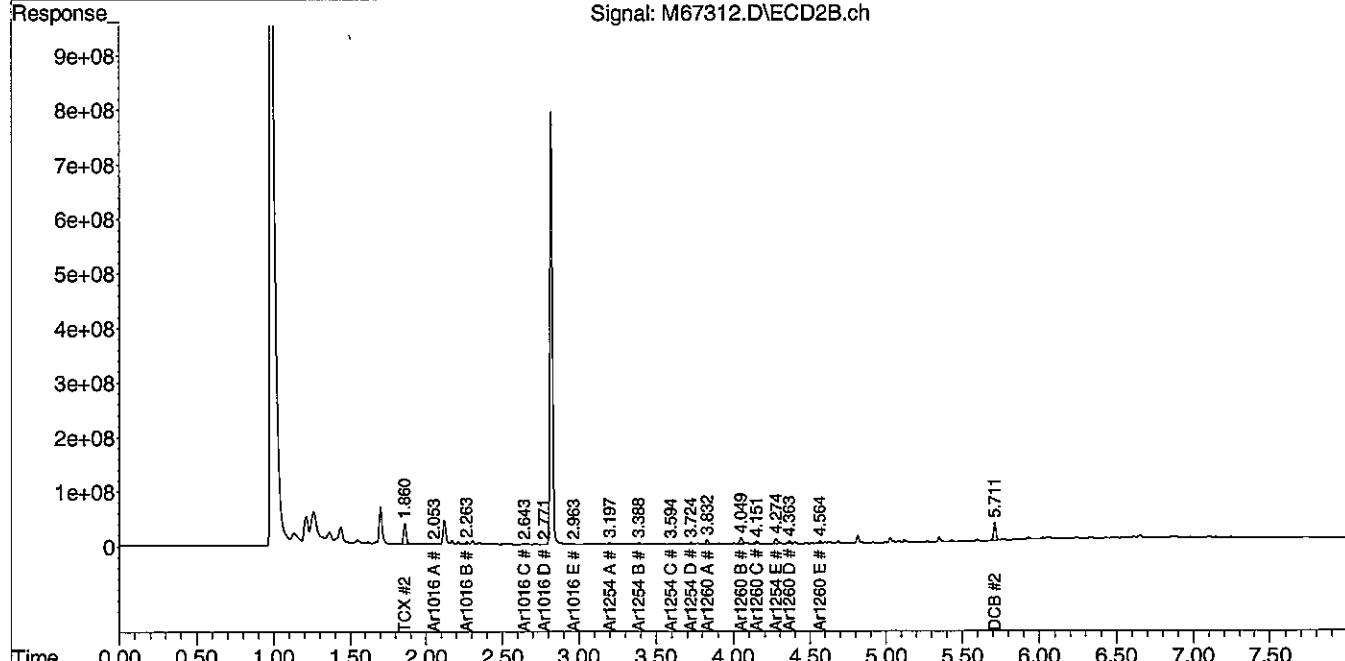
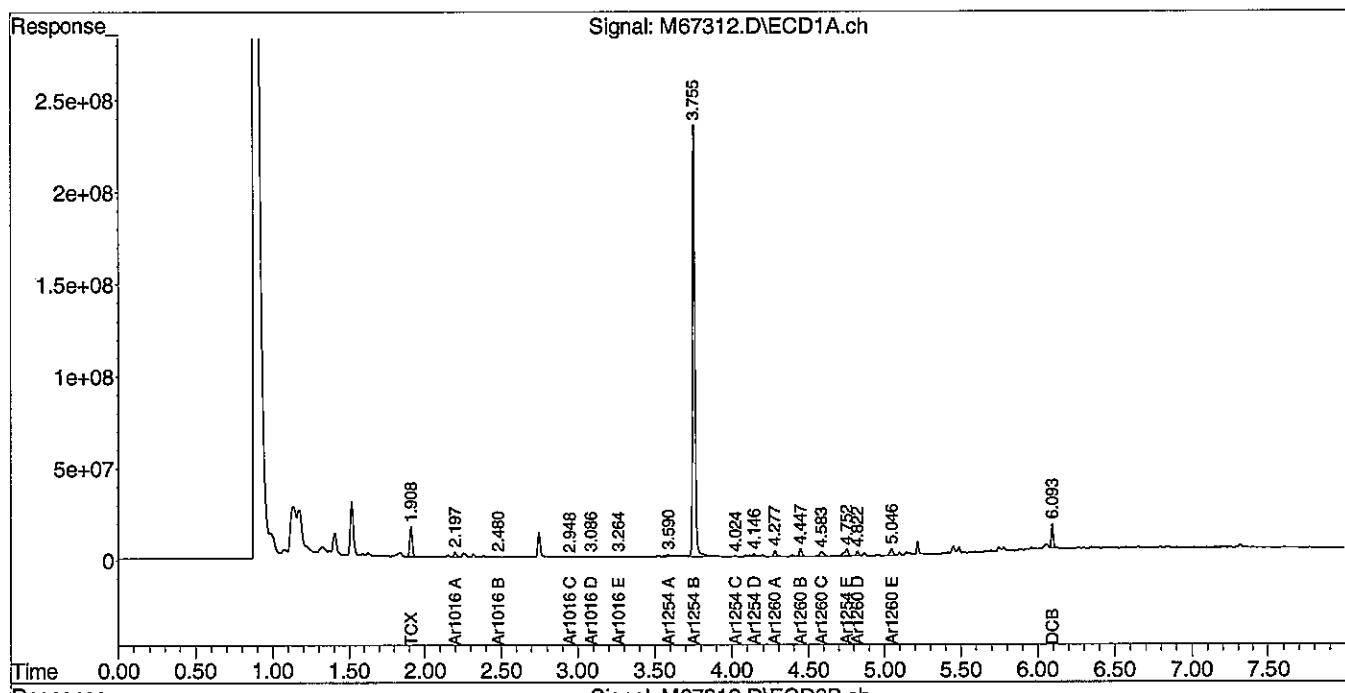
* Values outside QC limits

Comments: _____

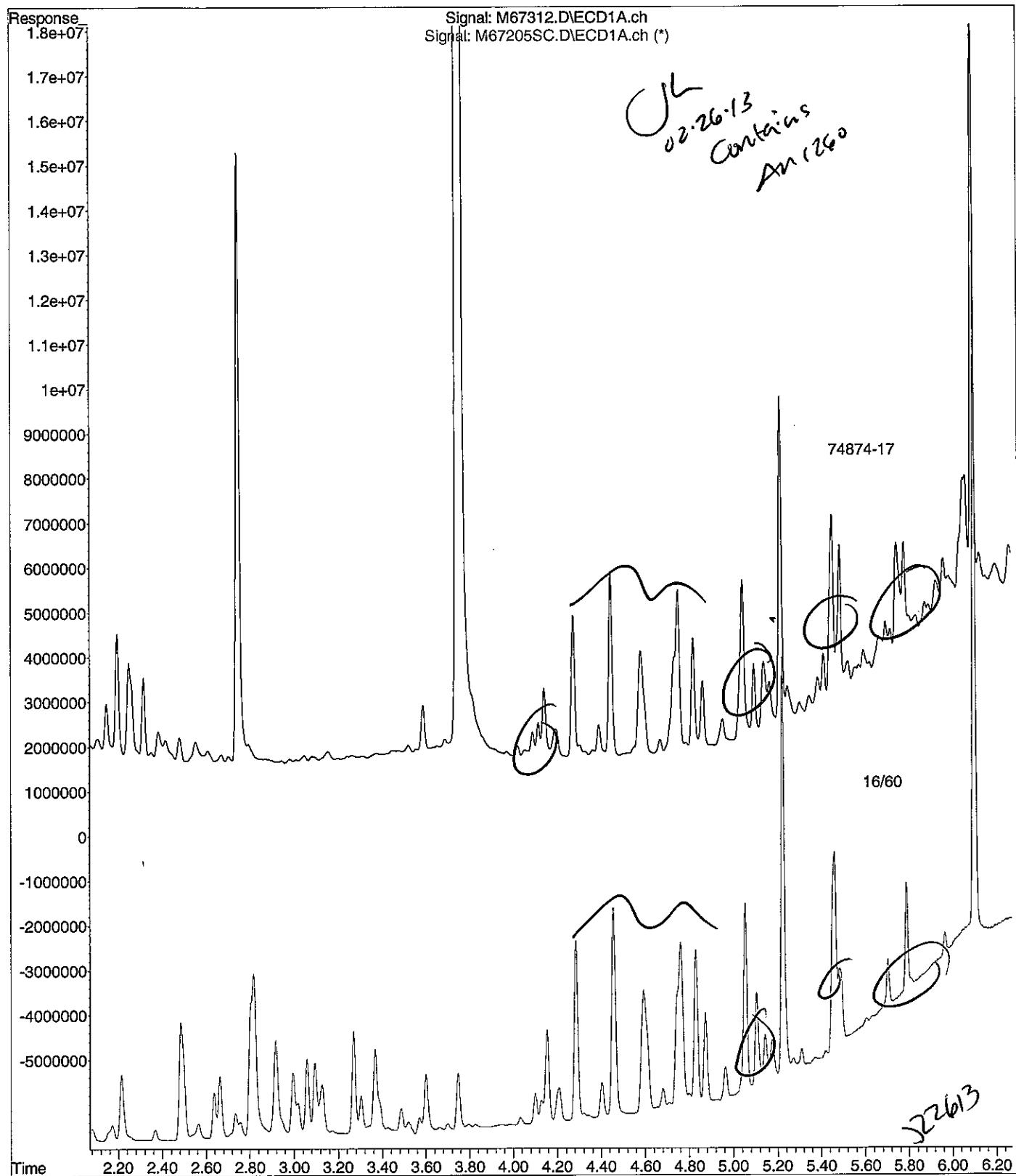
Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67312.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 23 Feb 2013 7:37 am
 Operator : JK
 Sample : 74874-17,,A/C
 Misc : SOIL
 ALS Vial : 48 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 16:01:28 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022213-M\M67312.D
Operator : JK
Acquired : 23 Feb 2013 7:37 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74874-17,,A/C
Misc Info : SOIL
Vial Number: 48



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-224 (6-8')

Lab Sample ID: 74874-18
Matrix: Solid
Percent Solid: 96
Dilution Factor: 20
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/21/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	660	U
PCB-1221	660	U
PCB-1232	660	U
PCB-1242	660	U
PCB-1248	660	U
PCB-1254	660	U
PCB-1260	660	9260

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M SDG: 74874
GC Column #1: STX-CLPesticides I Sample: 74874-18,1:20,,A/C
Column ID: 0.25 mm Data File: M67448.D
GC Column #2: STX-CLPesticides II Dilution Factor: 19.6
Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	9265	8097	13.5	

Column to be used to flag RPD values greater than QC limit of 40%

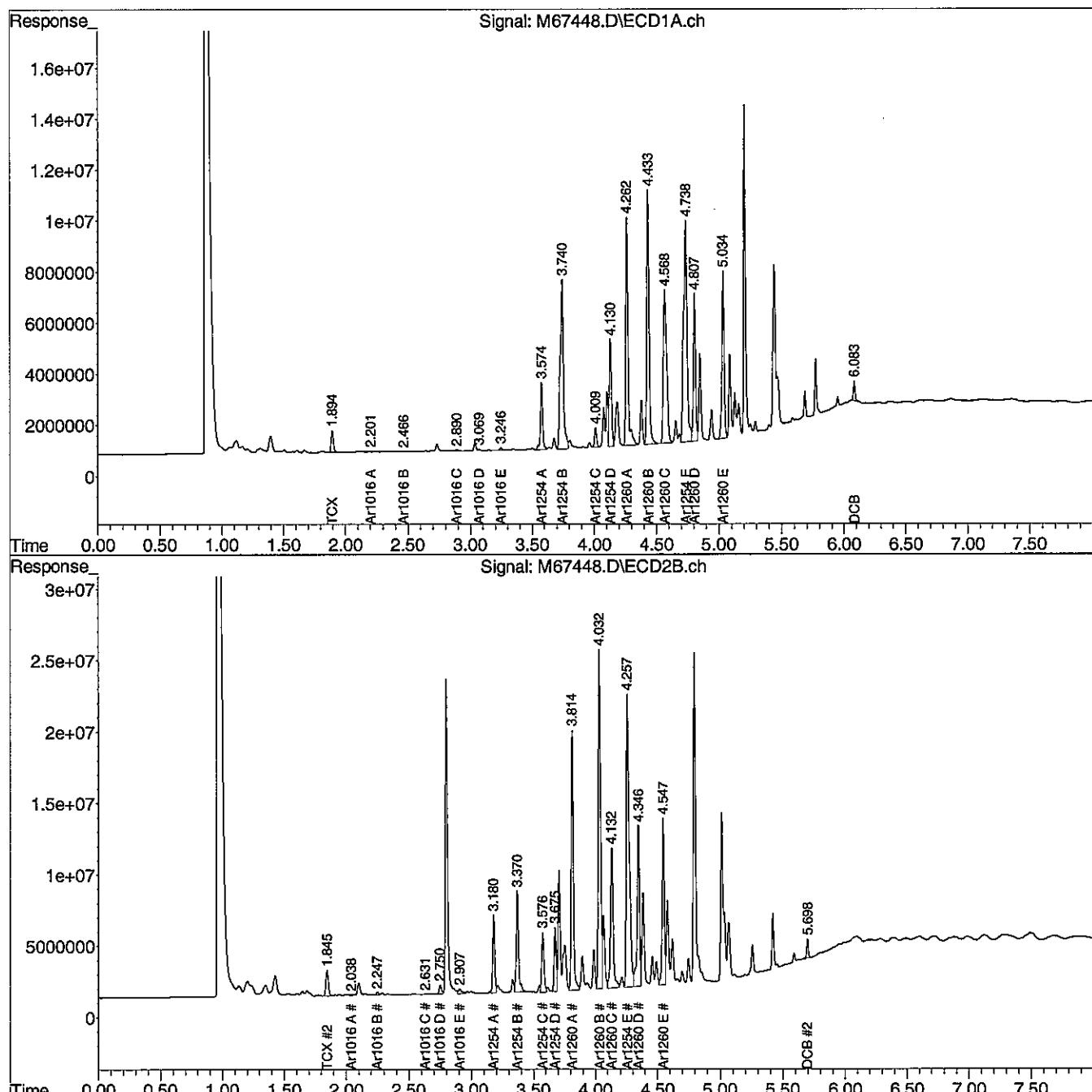
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67448.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:45 am
 Operator : JK
 Sample : 74874-18,1:20,,A/C
 Misc : SOIL
 ALS Vial : 36 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:58 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 26, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-225 (0-2')

Lab Sample ID:	74874-19
Matrix:	Solid
Percent Solid:	86
Dilution Factor:	1.1
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/23/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	36	U
PCB-1221	36	U
PCB-1232	36	U
PCB-1242	36	U
PCB-1248	36	U
PCB-1254	36	U
PCB-1260	36	310

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	68	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-19,,A/C

Column ID: 0.25 mm

Data File: M67314.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	297	310	4.0	

Column to be used to flag RPD values greater than QC limit of 40%

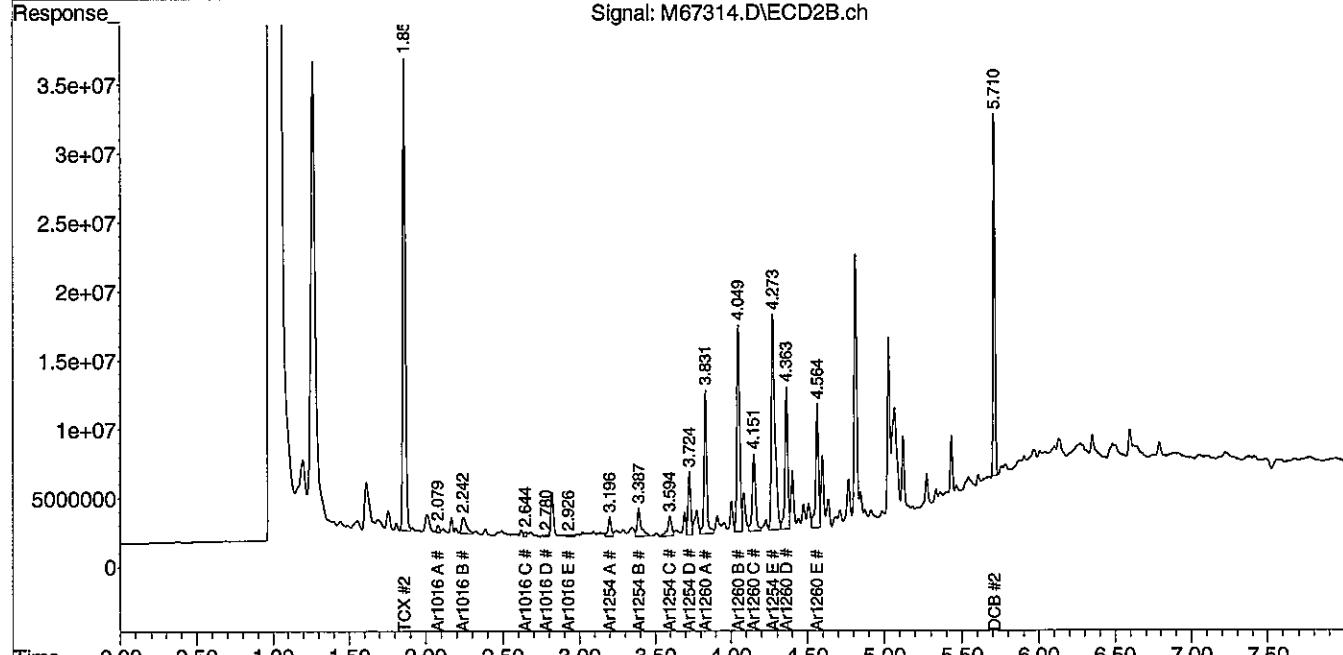
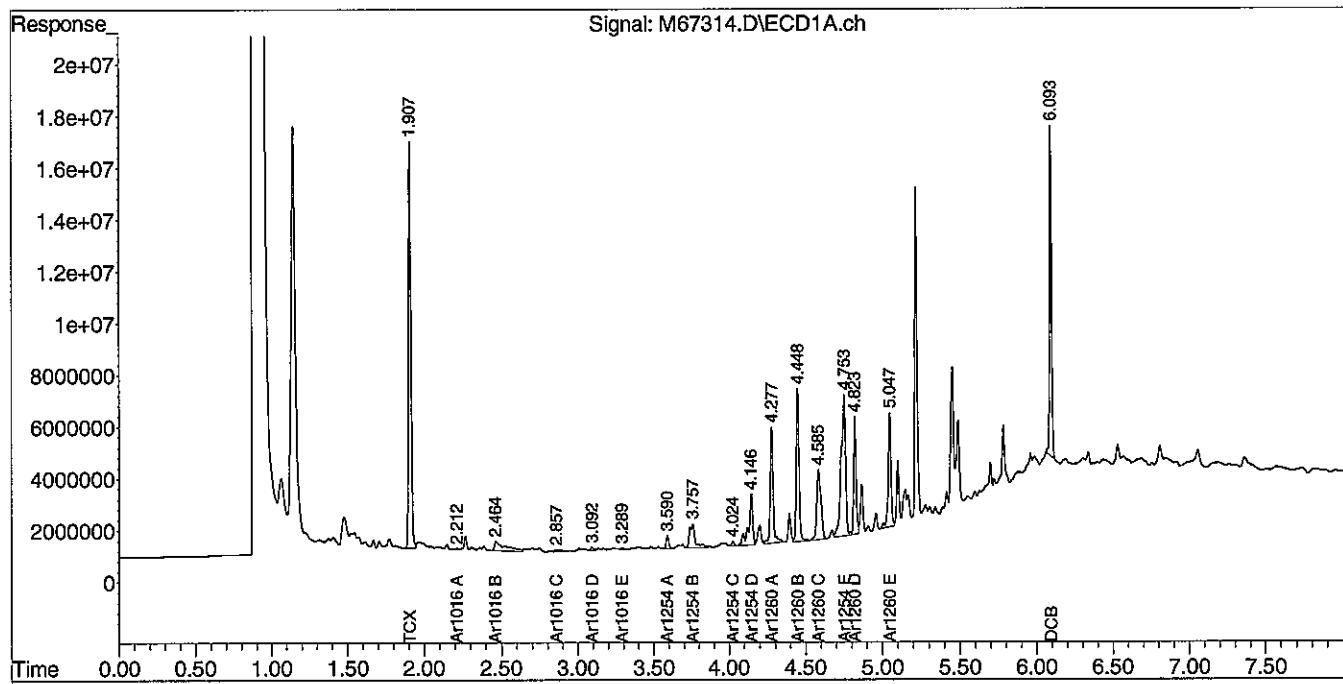
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67314.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 23 Feb 2013 7:57 am
 Operator : JK
 Sample : 74874-19,, A/C
 Misc : SOIL
 ALS Vial : 50 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 15:56:03 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:30 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-225 (2-4')

Lab Sample ID:	74874-20
Matrix:	Solid
Percent Solid:	90
Dilution Factor:	27
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/21/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	891	U
PCB-1221	891	U
PCB-1232	891	U
PCB-1242	891	U
PCB-1248	891	U
PCB-1254	891	U
PCB-1260	891	15800

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74874

GC Column #1: STX-CLPesticides I

Sample: 74874-20,1:25,,A/C

Column ID: 0.25 mm

Data File: M67449.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 27.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	15750	13908	12.4	

Column to be used to flag RPD values greater than QC limit of 40%

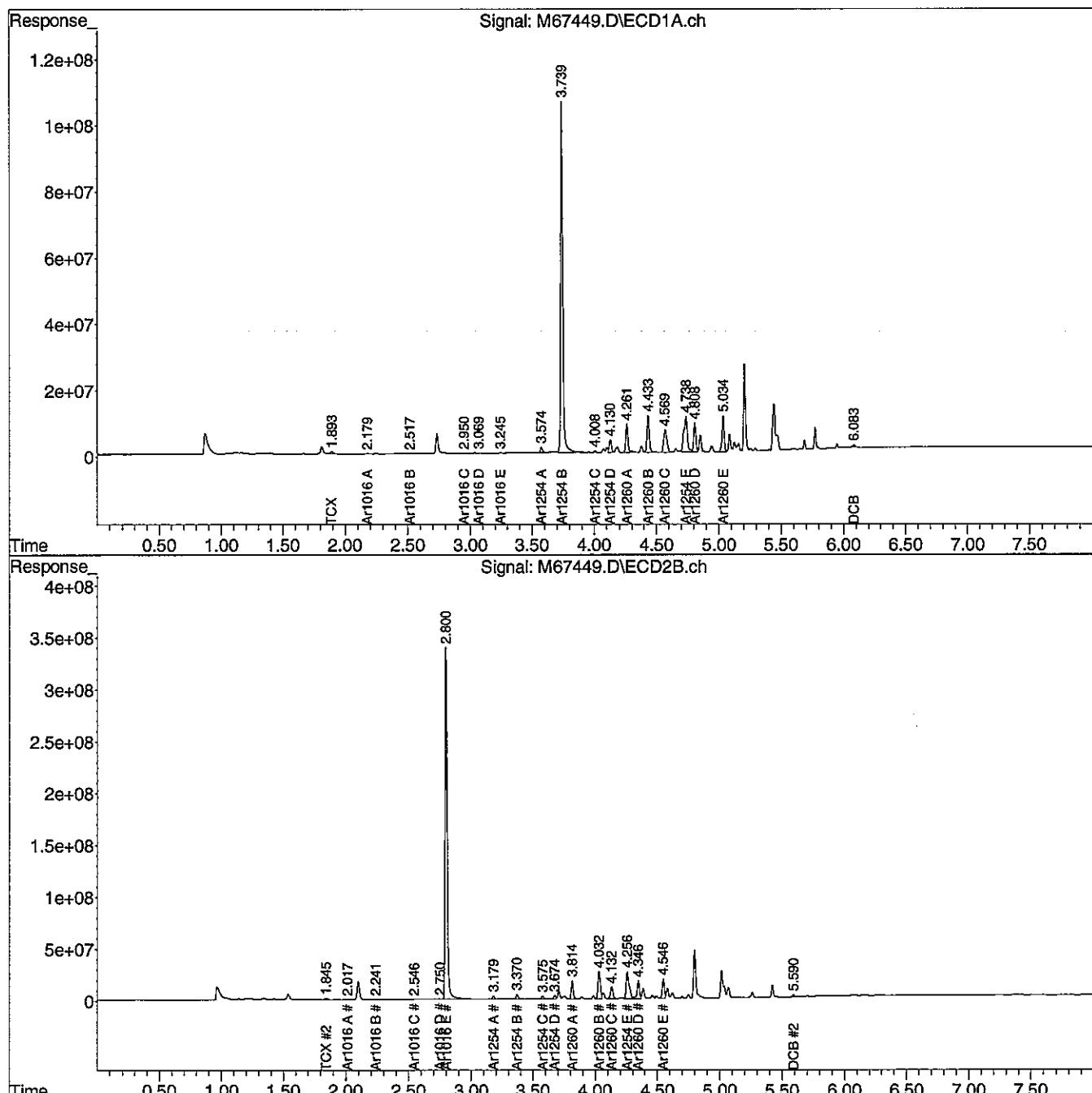
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67449.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 12:55 am
 Operator : JK
 Sample : 74874-20,1:25,,A/C
 Misc : SOIL
 ALS Vial : 37 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:48:00 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. :: 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



**PCB
QC FORMS**

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February 26, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	Lab QC

Lab Sample ID:	B022113PSOX RR
Matrix:	Soil
Percent Solid:	100
Dilution Factor:	1.0
Collection Date:	
Lab Receipt Date:	
Extraction Date:	02/21/13
Analysis Date:	02/23/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	112	%
Decachlorobiphenyl	113	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

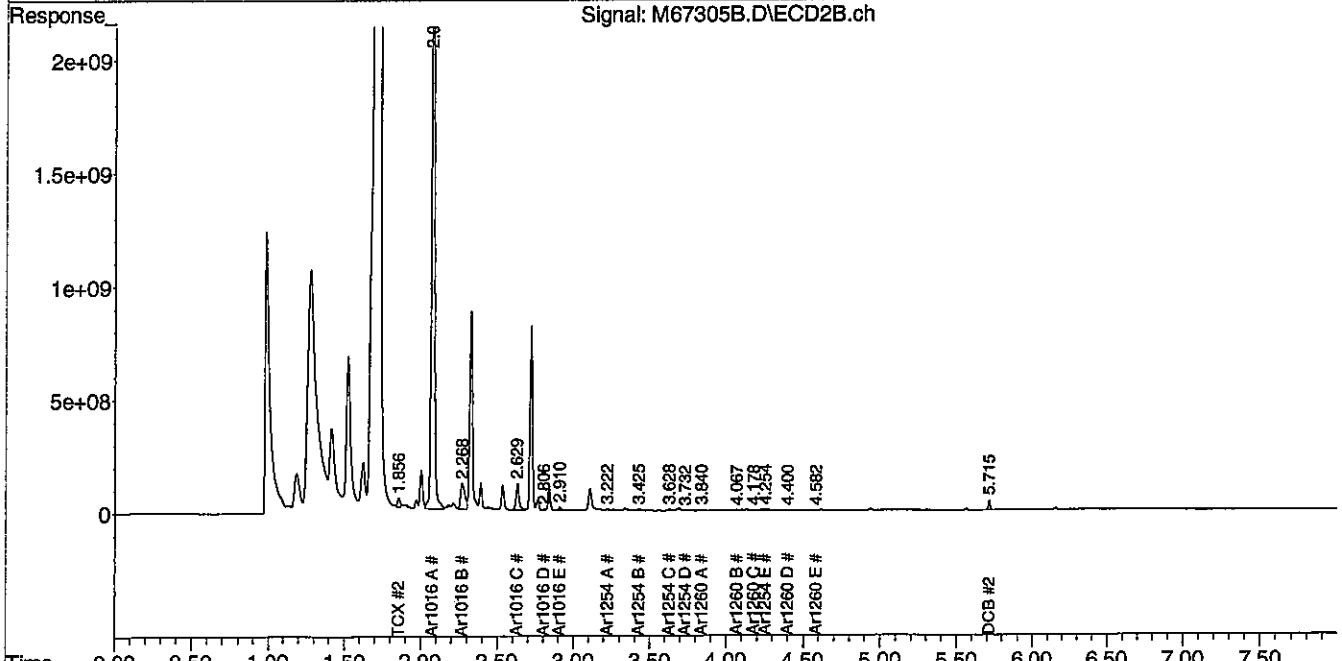
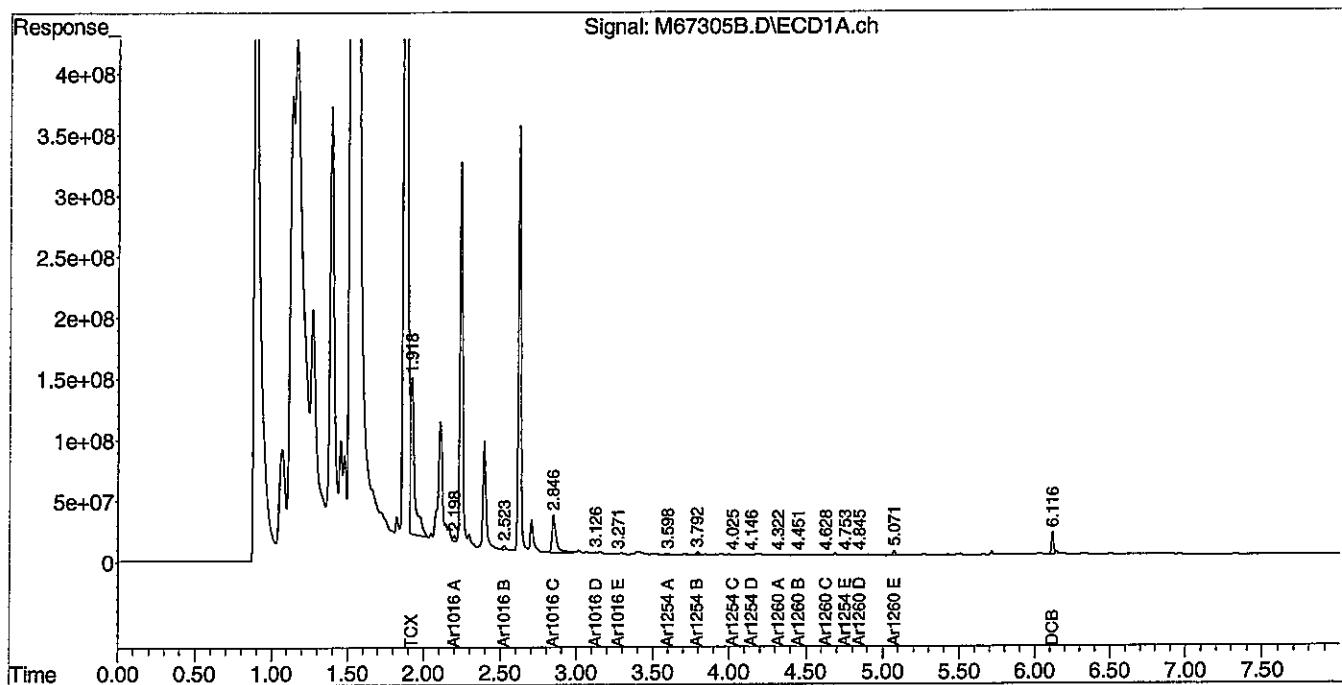
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022213-M\
 Data File : M67305B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 23 Feb 2013 6:26 am
 Operator : JK
 Sample : B022113PSOX,RR,,A/C
 Misc : SOIL
 ALS Vial : 28 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 26 15:56:46 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Feb 22 09:21:29 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022113PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/21/13
Analysis Date: 02/26/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	119	%
Decachlorobiphenyl	104	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

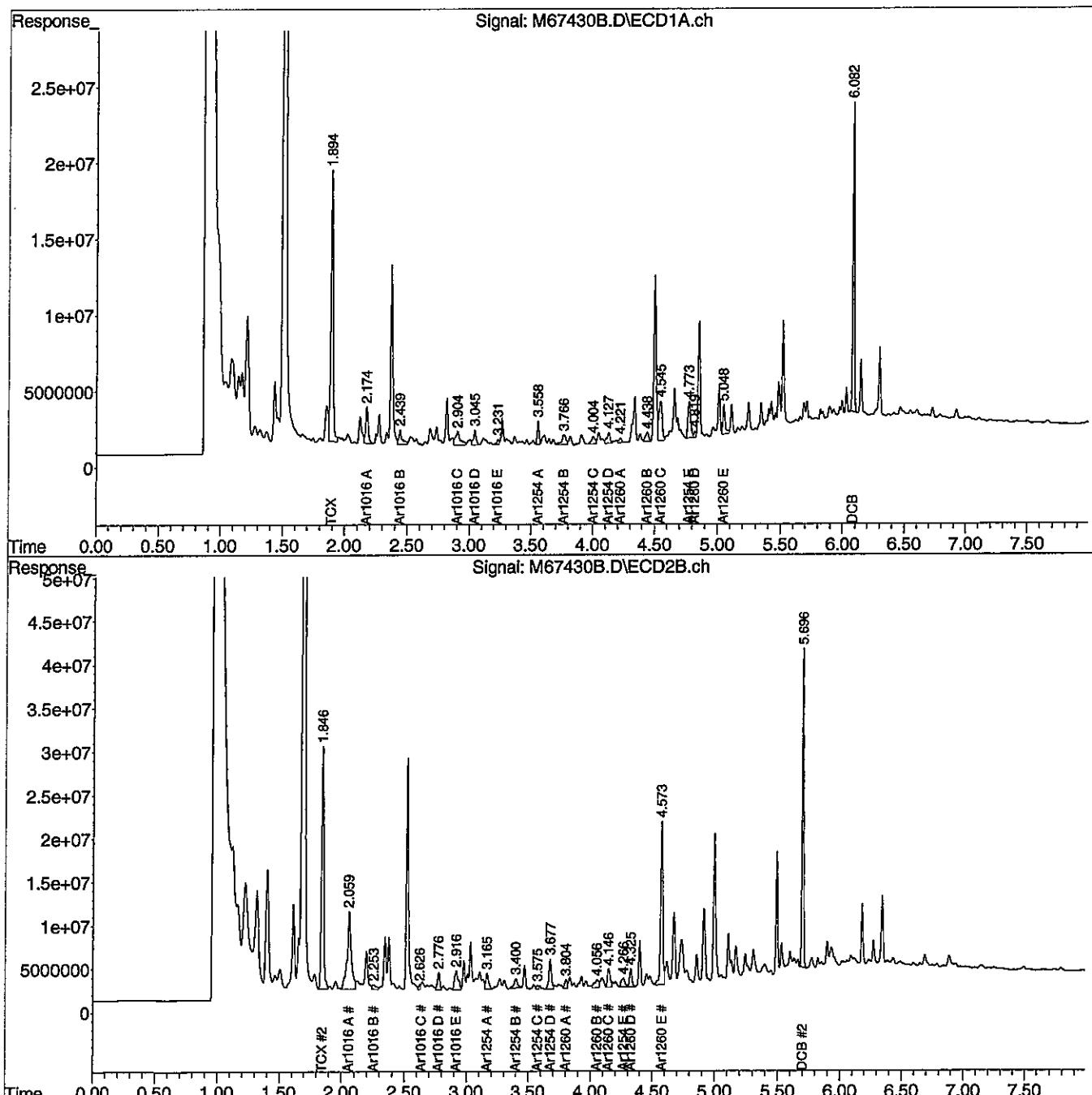
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67430B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 26 Feb 2013 9:44 pm
 Operator : JK
 Sample : B022113PSOX,,A/C
 Misc : SOIL
 ALS Vial : 18 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 11:47:22 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74874

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74874

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits
* Values outside QC limits
D System Monitoring Compound diluted out

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022113PSOX,,A/C

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: L022113PSOX,,A/C

Spike duplicate: LD022113PSOX,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP		
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC #	RESULT (ug/kg)	% REC #	RPD	#
PCB 1016	200	200	65	140	30	0	230	115	242	121		4.9
PCB 1260	200	200	60	130	30	0	211	105	239	119		12.6
PCB 1016 #2	200	200	65	140	30	0	190	95	229	115		19.0
PCB 1260 #2	200	200	60	130	30	0	171	85	193	97		12.4

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

**PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: 74874-2,,A/C

Spike: 74874-2,MS,,A/C

Spike duplicate: 74874-2,MSD,,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC #	RESULT (ug/kg)	% REC #	RPD #
PCB 1016	223	211	65	140	30	0	242	108	249	118	3.0
PCB 1260	223	211	60	130	30	308	439	59 *	447	65	1.7
PCB 1016 #2	223	211	65	140	30	0	193	87	190	90	1.7
PCB 1260 #2	223	211	60	130	30	268	372	46 *	388	57 *	4.3

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form

 Maine Energy Laboratory LLC 195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax																																																																																																																																																													
<p>Project Name: <u>MAINE ENERGY</u></p> <p>Project#: <u>12-3259.1</u></p> <p>Company: <u>Sunrise St.</u></p> <p>Report to: <u>John Cresssey</u></p> <p>Address: <u>164 Main St.</u></p> <p>Phone: <u>(207) 795-6009</u></p> <p>Quote #: <u></u></p> <p>PO# (if required): <u>12-3359.1</u></p> <p>Sample Identification</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th>Sample Date</th> <th>Sample Time</th> <th>Field Filtered by or N</th> <th>VOC: 8260 5242 624</th> <th>SVC: 8270 625 PAH only SIM</th> <th>Pestidles: 8081 608</th> <th>TPH: 8015 (Gas Range) ME4217</th> <th>TPH: 8015 (Diesel Range) ME4217</th> <th>EPH: Full or Ranges only TEPH</th> <th>VPH: Full or Ranges only</th> <th>Matrix Key:</th> </tr> </thead> <tbody> <tr> <td>2-19</td> <td>1432</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>A= HCl B= 4 C C= Unpres D = HNO3 E = H2SO4 F = Hexane G = Other</td> </tr> <tr> <td>SE-58-221 (2-4)</td> <td>1432</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>B= 4 C</td> </tr> <tr> <td>SE-58-221 (4-5)</td> <td>1441</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>C= Unpres</td> </tr> <tr> <td>SE-58-220 (0-2)</td> <td>1515</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>D = HNO3</td> </tr> <tr> <td>SE-58-220 (2-4)</td> <td>1524</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>E = H2SO4</td> </tr> <tr> <td>SE-58-220 (4-6)</td> <td>1530</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>F = Hexane</td> </tr> <tr> <td>SE-58-220 (6-8)</td> <td>1535</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>G = Other</td> </tr> <tr> <td>SE-58-220 (8-10)</td> <td>1545</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> <tr> <td>SE-58-220 (10-11)</td> <td>1550</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> <tr> <td>SE-58-220 (0-2)</td> <td>1557</td> <td>N</td> <td>X</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> </tr> </tbody> </table> <p>Preservation Code:</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th>Preservation Key:</th> <th>1) Shipped or hand delivered</th> <th>2) Temperature (°C):</th> <th>3) Received in good condition:</th> <th>4) pH checked by:</th> <th>5) Labels checked by:</th> </tr> </thead> <tbody> <tr> <td>A= HCl B= 4 C C= Unpres D = HNO3 E = H2SO4 F = Hexane G = Other</td> <td><input checked="" type="checkbox"/></td> </tr> </tbody> </table> <p>Comments, Additional Analyses, or Special Instructions:</p> <p>Email Results to: <u>John Cresssey</u></p> <p>Turnaround Time (TAT)</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th>1 Day*</th> <th>2 Days*</th> <th>4 Days*</th> <th>Standard</th> </tr> </thead> <tbody> <tr> <td><input type="checkbox"/></td> <td><input type="checkbox"/></td> <td><input type="checkbox"/></td> <td><input checked="" type="checkbox"/></td> </tr> </tbody> </table> <p>*Fee may apply; lab approval required</p> <p>Missing sample SE-58-21(0-2) @ 15:57 ↳ No sample collected per client. 02/20/13 -kg</p> <p>Project Requirements:</p> <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th>Report Type:</th> <th>State:</th> <th>State Standard:</th> </tr> </thead> <tbody> <tr> <td><input type="checkbox"/> MCP*</td> <td><input type="checkbox"/> Level II*</td> <td><input type="checkbox"/> NH</td> </tr> <tr> <td><input type="checkbox"/> CTRCP*</td> <td><input type="checkbox"/> Level III*</td> <td><input type="checkbox"/> MA</td> </tr> <tr> <td><input type="checkbox"/> DOD*</td> <td><input type="checkbox"/> Level IV*</td> <td><input checked="" type="checkbox"/> ME <input type="checkbox"/> CT <input type="checkbox"/> RI</td> </tr> <tr> <td><input checked="" type="checkbox"/> Standard</td> <td><input type="checkbox"/> Other:</td> <td><input type="checkbox"/> Other: _____ <i>[Signature]</i></td> </tr> </tbody> </table> <p>EDD Required: <u>Y/N</u></p> <p>Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.</p> <p>Relinquished By Sampler: <u>John Cresssey</u></p> <p>Date: <u>2-20</u> Time: <u>1731</u> Received By: <u>John Cresssey</u></p> <p>Relinquished By: <u></u></p> <p>Date: <u></u> Time: <u></u> Received By: <u></u></p> <p>Relinquished By: <u></u></p> <p>Date: <u></u> Time: <u></u> Received By: <u></u></p>				Sample Date	Sample Time	Field Filtered by or N	VOC: 8260 5242 624	SVC: 8270 625 PAH only SIM	Pestidles: 8081 608	TPH: 8015 (Gas Range) ME4217	TPH: 8015 (Diesel Range) ME4217	EPH: Full or Ranges only TEPH	VPH: Full or Ranges only	Matrix Key:	2-19	1432	N	X							A= HCl B= 4 C C= Unpres D = HNO3 E = H2SO4 F = Hexane G = Other	SE-58-221 (2-4)	1432	N	X							B= 4 C	SE-58-221 (4-5)	1441	N	X							C= Unpres	SE-58-220 (0-2)	1515	N	X							D = HNO3	SE-58-220 (2-4)	1524	N	X							E = H2SO4	SE-58-220 (4-6)	1530	N	X							F = Hexane	SE-58-220 (6-8)	1535	N	X							G = Other	SE-58-220 (8-10)	1545	N	X								SE-58-220 (10-11)	1550	N	X								SE-58-220 (0-2)	1557	N	X								Preservation Key:	1) Shipped or hand delivered	2) Temperature (°C):	3) Received in good condition:	4) pH checked by:	5) Labels checked by:	A= HCl B= 4 C C= Unpres D = HNO3 E = H2SO4 F = Hexane G = Other	<input checked="" type="checkbox"/>	1 Day*	2 Days*	4 Days*	Standard	<input type="checkbox"/>	<input type="checkbox"/>	<input type="checkbox"/>	<input checked="" type="checkbox"/>	Report Type:	State:	State Standard:	<input type="checkbox"/> MCP*	<input type="checkbox"/> Level II*	<input type="checkbox"/> NH	<input type="checkbox"/> CTRCP*	<input type="checkbox"/> Level III*	<input type="checkbox"/> MA	<input type="checkbox"/> DOD*	<input type="checkbox"/> Level IV*	<input checked="" type="checkbox"/> ME <input type="checkbox"/> CT <input type="checkbox"/> RI	<input checked="" type="checkbox"/> Standard	<input type="checkbox"/> Other:	<input type="checkbox"/> Other: _____ <i>[Signature]</i>				
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Chain Of Custody Form

 environmental laboratory LLC		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 928-9906		(603) 436-5111 (603) 430-2151 Fax	For Analytics Use Only		
Project Name: MAIN ENERGY Project #: 12-3259.1		Preservation Code: Please <input checked="" type="checkbox"/> in preservation code here		Circle and/or Write Required Analysis Followed by Preservation Code <input checked="" type="checkbox"/> * * *		Samples were: 0-3-0-7c	
Company:	Summit	Preservation Key: A = HCL B = 4°C C = Unpres D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other	Sample Date	Sample Time	Matrix	No. of Containers checked	Analytics Sample #
Report to:	John Cressy	2-19	1402	S	1	-11	
Address:	640 MAIN ST. Lewiston ME 04240	2-19	1655	S/0	1	-12	
Phone:	(207) 795-6069	2-19	1320	X	1	-13	
Quote #:		2-19	1622	X	1	-14	
PO# (if required):	12-33-59.1	2-19	1638	X	1	-15	
		2-19	1650	X	1	-16	
		2-19	1700	X	1	-17	
		2-19	1734	X	1	-18	
		2-19	1736	X	1	-19	
						-20	
Comments, Additional Analyses, or Special Instructions:							
** List requested metals here <input checked="" type="checkbox"/> Fee may apply							
Email Results to: JCressy@summitbi.com		Report Type: <input type="checkbox"/> MCP* <input type="checkbox"/> Level II* <input type="checkbox"/> CTRCP* <input type="checkbox"/> Level III* <input type="checkbox"/> DOD* <input type="checkbox"/> Level IV* <input checked="" type="checkbox"/> Standard <input checked="" type="checkbox"/> Other		State Standard: <input type="checkbox"/> NH <input type="checkbox"/> MA <input checked="" type="checkbox"/> ME <input type="checkbox"/> CT <input type="checkbox"/> RI EDD Required: Y= N			
Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> 4 Days* <input type="checkbox"/> 5 Days <input checked="" type="checkbox"/> Standard <small>(6-10 business days)</small>							
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.							
Relinquished By Sampler: JCressy		Date: 3-20	Time: 11:13	Received By: LC Stevens			
Relinquished By: 		Date:	Time:	Received By:			
Relinquished By: 		Date:	Time:	Received By:			

*Fee may apply; lab approval required



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 74874
CLIENT: Summit
PROJECT: Main Energy

COOLER NUMBER: 74,223
NUMBER OF COOLERS: 2

A: PRELIMINARY EXAMINATION:

1. Cooler received by (initials): KB

DATE COOLER RECEIVED/OPENED: 02/20/13

2. Circle one:

Hand delivered
(If no, skip 3)

Shipped

3. Did cooler come with a shipping slip?

Y

N

3a. Enter carrier name and airbill number here:

4. Were custody seals on the outside of cooler?

How many & where:

Seal Date:

Y

N

Seal Name:

5. Did the custody seals arrive unbroken and intact upon arrival?

Y

N

6. COC#:

7. Were Custody papers filled out properly (ink, signed, legible, project information etc)?

Y

N

8. Were custody papers sealed in a plastic bag?

Y

N

9. Did you sign the COC in the appropriate place?

Y

N

10. Was enough ice used to chill the cooler?

Y

N Temp. of cooler:

0.3°C, 0.7°C

B. Log-In: Date samples were logged in:

KB By: 02/20/13

11. Were all bottles sealed in separate plastic bags?

Y

N

12. Did all bottles arrive unbroken and were labels in good condition?

Y

N

13. Were all bottle labels complete (ID, Date, time, etc.)

Y

N

14. Did all bottle labels agree with custody papers?

Y

N

15. Were the correct containers used for the tests indicated?

Y

N

16. Were samples received at the correct pH?

Y

N

17. Was sufficient amount of sample sent for the tests indicated?

Y

N

18. Were all samples submitted within holding time?

Y

N

19. Were all containers used within AEL's expiration date?**

Y

N

20. Were VOA samples absent of greater than pea-sized bubbles?

Y

N

(Note: Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

*If NO, List Sample ID's, Lab #s:

When bubbles are present in VOA samples they are labelled from smallest (or no bubbles) to largest. Lab to analyze VOA samples with no bubbles or smallest bubbles first

20. Laboratory labeling verified by (initials): KB

Date: 2/20/13

**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.

February 27, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74875
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples
PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
Chromatograms
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No exceptions.

PCBs by EPA Method 8082:
No results were reported below the quantitation limit.

All samples except 74875-10 and 74875-12 thru 74875-14 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument.

The MS/MSD analyzed on sample 74875-7 was only assessed for precision of PCB 1260 due to the parent sample having concentrations of PCB 1260 that exceeded the calibration range of the instrument. The RPDs for the MS/MSD did not meet acceptance criteria. The laboratory control samples (L022213PSOX/LD022213PSOX) were in control for all recoveries and RPDs. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC



Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
Summit Environmental Consultants Inc.
640 Main Street
Lewiston ME 04240

Report Number: 74875

Revision: Rev. 0

Re: MAINE ENERGY (Project No: 12-3259.1)

Enclosed are the results of the analyses on your sample(s). Samples were received on 20 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Sample Analysis: The attached pages detail the Client Sample IDs, Lab Sample IDs, and Analyses requested

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

If you have any questions on these results, please do not hesitate to contact us.

Authorized signature



Stephen L. Knollmeyer Lab. Director

Date

02/28/2013

**This report shall not be reproduced, except in full, without the written
consent of Analytics Environmental Laboratory, LLC.**

CLIENT: Summit Environmental
 Consultants Inc.

REPORT NUMBER: 74875

REV: Rev. 0

PROJECT: MAINE ENERGY (Project No: 12-3259.1)

<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74875-1	02/19/13	SE-SB-225 (4-6')	EPA 8082 (PCBs only)	
74875-2	02/19/13	SE-SB-225 (6-7.5')	EPA 8082 (PCBs only)	
74875-3	02/20/13	SE-SB-226 (0-2')	EPA 8082 (PCBs only)	
74875-4	02/20/13	SE-SB-226 (2-4')	EPA 8082 (PCBs only)	
74875-5	02/20/13	SE-SB-226 (4-5')	EPA 8082 (PCBs only)	
74875-6	02/20/13	SE-SB-232 (0-2')	EPA 8082 (PCBs only)	
74875-7	02/20/13	SE-SB-232A (2-4')	EPA 8082 (PCBs only)	
74875-8	02/20/13	SE-SB-232A (4-6')	EPA 8082 (PCBs only)	
74875-9	02/20/13	SE-SB-232A (6-8')	EPA 8082 (PCBs only)	
74875-10	02/20/13	SE-SB-232A (8-10')	EPA 8082 (PCBs only)	
74875-11	02/20/13	SE-SB-233 (0-2')	EPA 8082 (PCBs only)	
74875-12	02/20/13	SE-SB-233 (2-4')	EPA 8082 (PCBs only)	
74875-13	02/20/13	SE-SB-233 (4-6')	EPA 8082 (PCBs only)	
74875-14	02/20/13	SE-SB-233 (6-8')	EPA 8082 (PCBs only)	
74875-15	02/20/13	SE-SB-234 (0-2')	EPA 8082 (PCBs only)	

Surrogate Compound Limits

	Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water				
1,4-Difluorobenzene		70-130		EPA 524.2
Bromofluorobenzene		70-130		
1,2-Dichlorobenzene-d4		70-130		
Volatile Organic Compounds				
1,2-Dichloroethane-d4		70-120	70-120	EPA 624/8260B
Toluene-d8		85-120	85-120	
Bromofluorobenzene		75-120	75-120	
Semi-Volatile Organic Compounds				
2-Fluorophenol		20-110	35-105	EPA 625/8270C
d5-Phenol		15-110	40-100	
d5-nitrobenzene		40-110	35-100	
2-Fluorobiphenyl		50-110	45-105	
2,4,6-Tribromophenol		40-110	40-125	
d14-p-terphenyl		50-130	30-125	
PAH's by SIM				
d5-nitrobenzene		21-110	35-110	EPA 8270C
2-Fluorobiphenyl		36-121	45-105	
d14-p-terphenyl		33-141	30-125	
Pesticides and PCBs				
2,4,5,6-Tetrachloro-m-xylene (TCX)		46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)		40-135	40-130	
Herbicides				
Dichloroacetic acid (DCAA)		30-150	30-150	
Gasoline Range Organics/TPH Gasoline				
Trifluorotoluene TFT (FID)		60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)		60-140	60-140	
Trifluorotoluene TFT (PID)		60-140	60-140	
Bromofluorobenzene (BFB) (PID)		60-140	60-140	
Diesel Range Organics/TPH Diesel				
m-terphenyl		60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons				
2,5-Dibromotoluene (PID)		70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)		70-130	70-130	
Extricable Petroleum Hydrocarbons				
1-chloro-octadecane (aliphatic)		40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)		40-140	40-140	
2-Fluorobiphenyl (Fractionation)		40-140	40-140	
2-Bromonaphthalene (fractionation)		40-140	40-140	

PCB
DATA SUMMARIES

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-225 (4-6')

Lab Sample ID:	74875-1
Matrix:	Solid
Percent Solid:	79
Dilution Factor:	24
Collection Date:	02/19/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/22/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	792	U
PCB-1221	792	U
PCB-1232	792	U
PCB-1242	792	U
PCB-1248	792	U
PCB-1254	792	U
PCB-1260	792	17600

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-1,1:20,,A/C

Column ID: 0.25 mm

Data File: M67505.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 24.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	17595	15559	12.3	

Column to be used to flag RPD values greater than QC limit of 40%

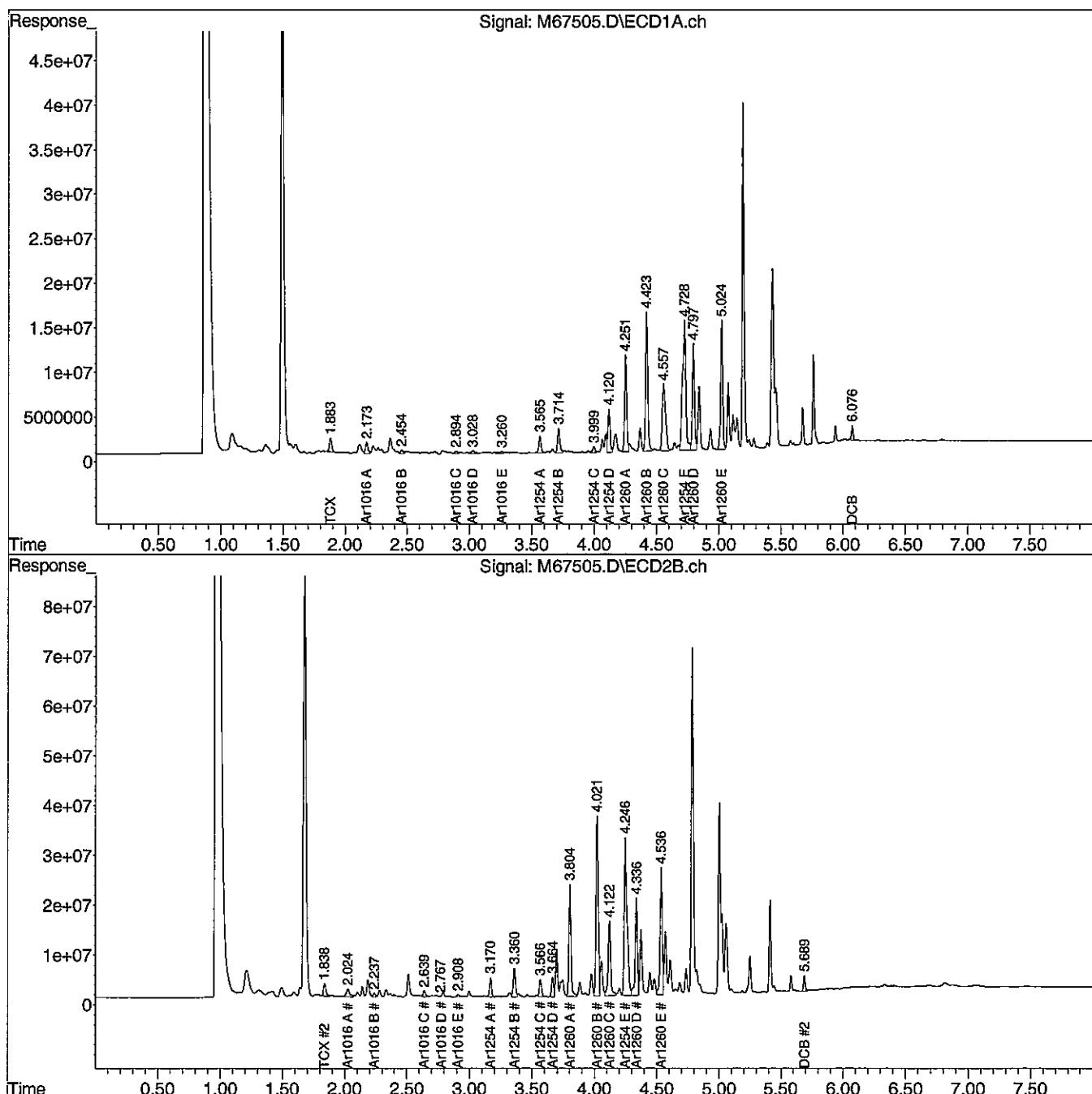
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67505.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:39 pm
 Operator : JK
 Sample : 74875-1,1:20,,A/C
 Misc : SOIL
 ALS Vial : 8 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:44:29 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-225 (6-7.5')

Lab Sample ID: 74875-2
Matrix: Solid
Percent Solid: 94
Dilution Factor: 21
Collection Date: 02/19/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	693	U
PCB-1221	693	U
PCB-1232	693	U
PCB-1242	693	U
PCB-1248	693	U
PCB-1254	693	U
PCB-1260	693	10900

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-2,1:20,,A/C

Column ID: 0.25 mm

Data File: M67506.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 21.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	10891	9752	11.0	

Column to be used to flag RPD values greater than QC limit of 40%

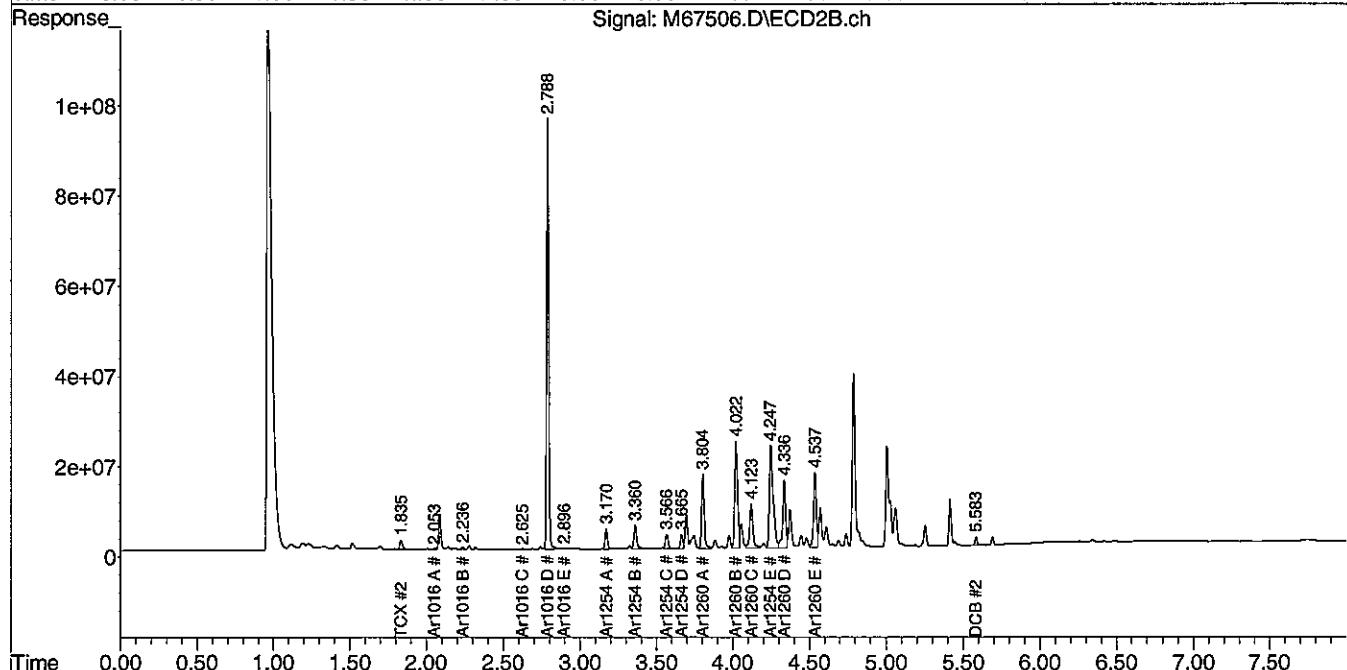
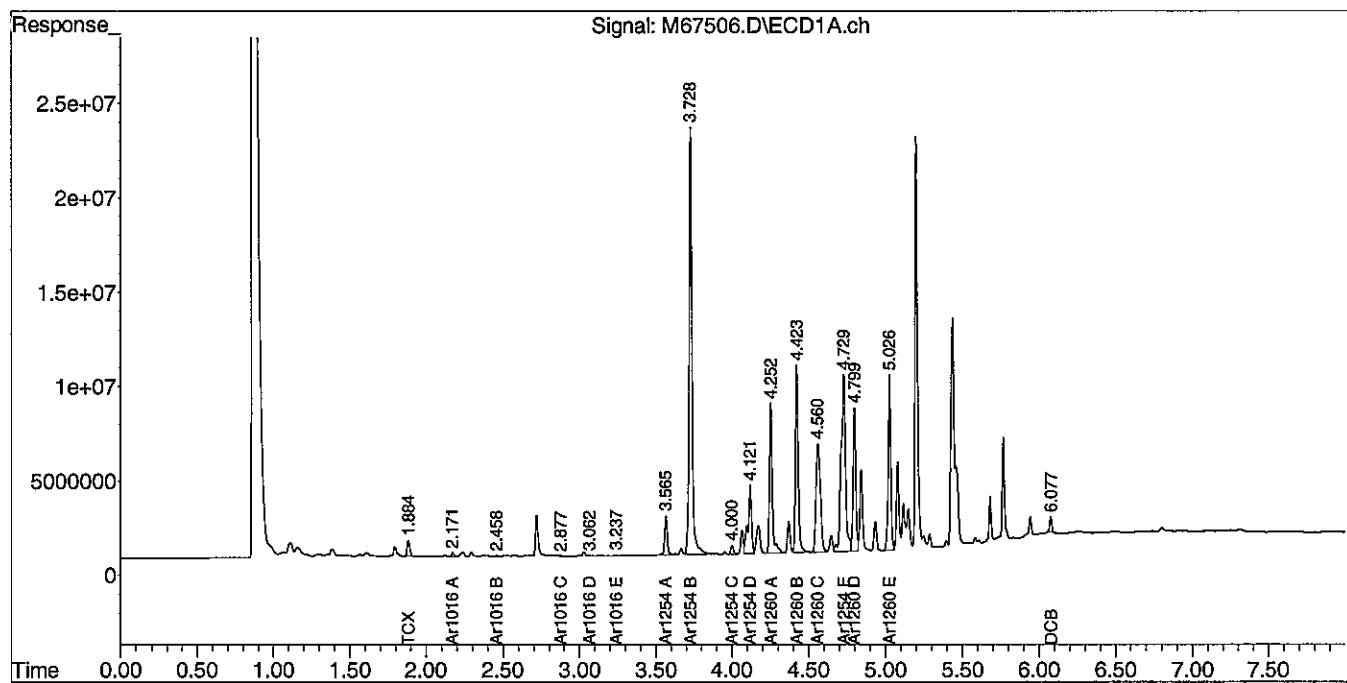
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67506.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:49 pm
 Operator : JK
 Sample : 74875-2,1:20,,A/C
 Misc : SOIL
 ALS Vial : 9 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:25 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-226 (0-2")

Lab Sample ID: 74875-3
 Matrix: Solid
 Percent Solid: 81
 Dilution Factor: 1230
 Collection Date: 02/20/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/22/13
 Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	40600	U
PCB-1221	40600	U
PCB-1232	40600	U
PCB-1242	40600	U
PCB-1248	40600	U
PCB-1254	40600	U
PCB-1260	40600	631000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-3,1:1000,,A/C

Column ID: 0.25 mm

Data File: M67526.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1228.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	631090	577563	8.9	

Column to be used to flag RPD values greater than QC limit of 40%

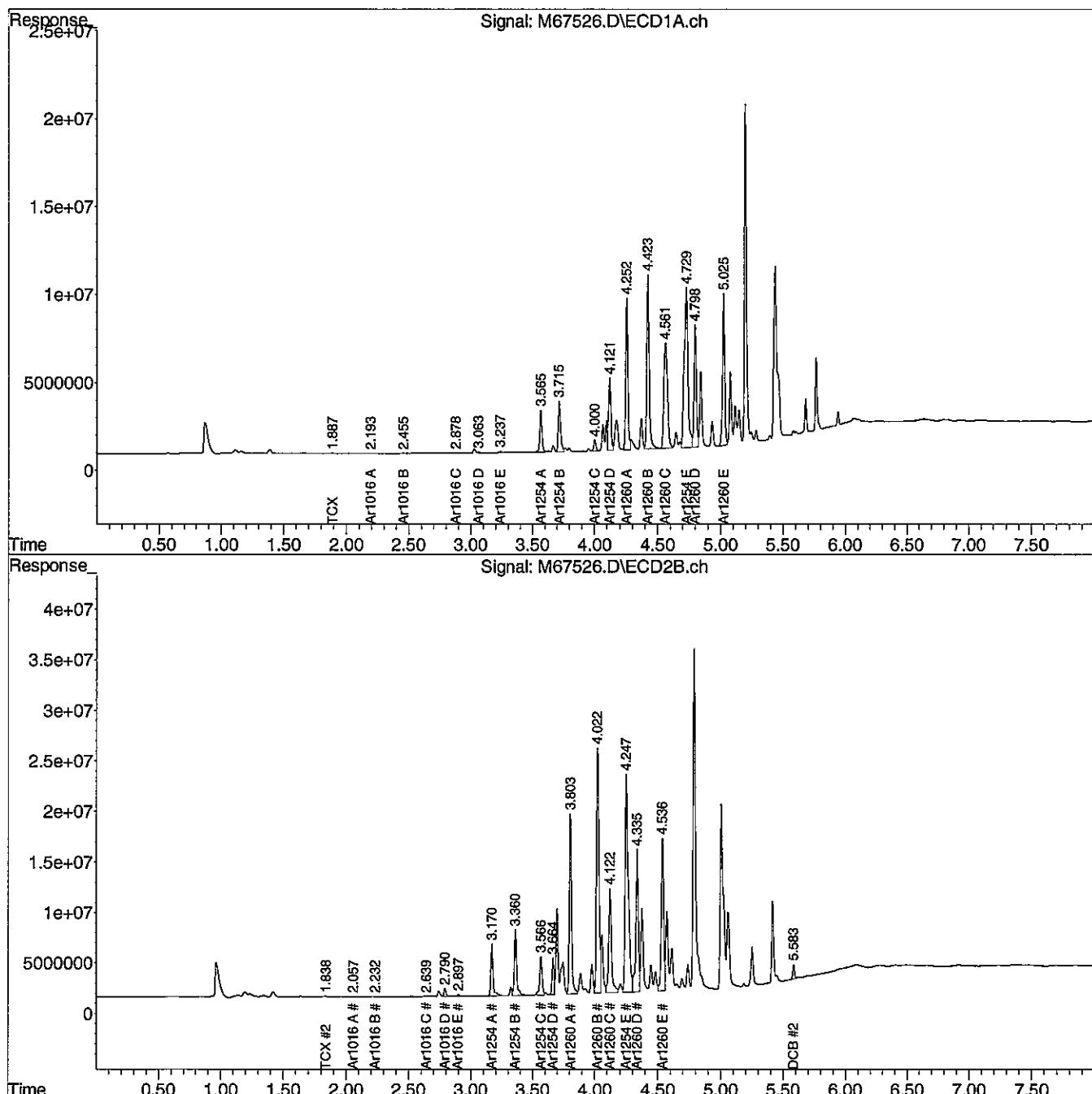
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67526.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 5:11 pm
 Operator : JK
 Sample : 74875-3,1:1000,,A/C
 Misc : SOIL
 ALS Vial : 7 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 17:31:03 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-226 (2-4")

Lab Sample ID:	74875-4
Matrix:	Solid
Percent Solid:	85
Dilution Factor:	23600
Collection Date:	02/20/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/22/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	779000	U
PCB-1221	779000	U
PCB-1232	779000	U
PCB-1242	779000	U
PCB-1248	779000	U
PCB-1254	779000	U
PCB-1260	779000	9660000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-4,1:20000,,A/C

Column ID: 0.25 mm

Data File: M67527.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 23589.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	9662352	8687004	10.6	

Column to be used to flag RPD values greater than QC limit of 40%

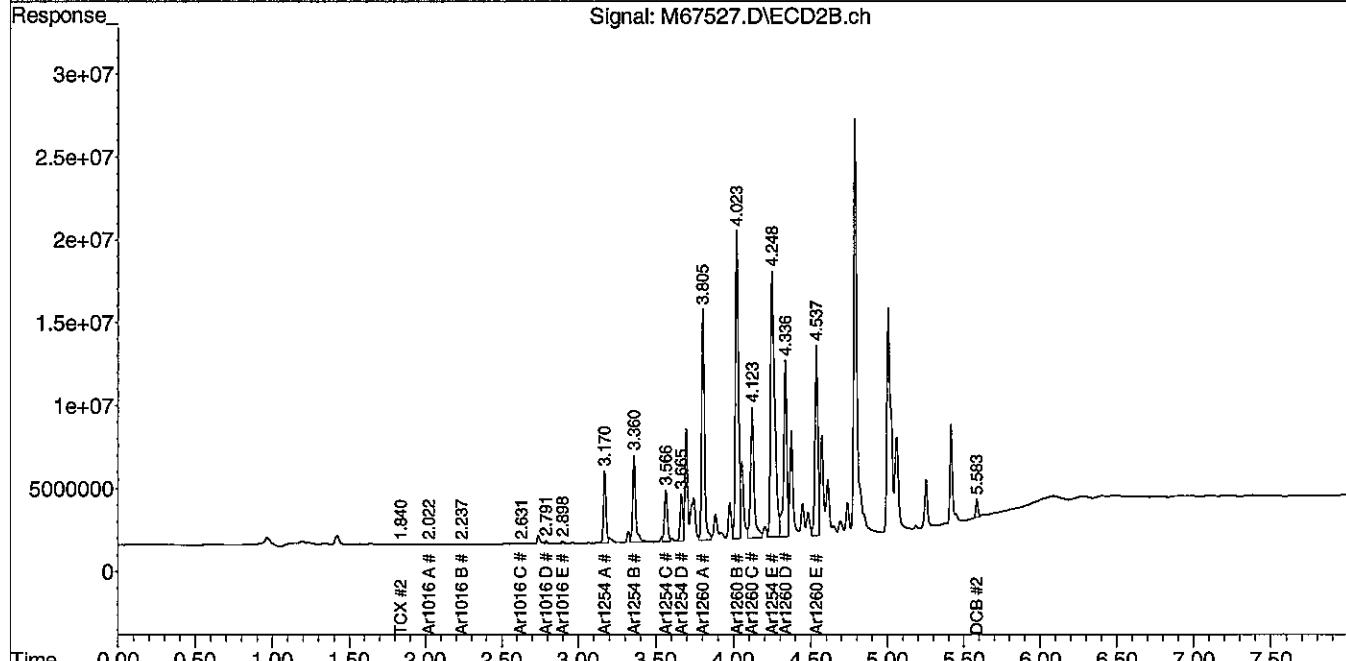
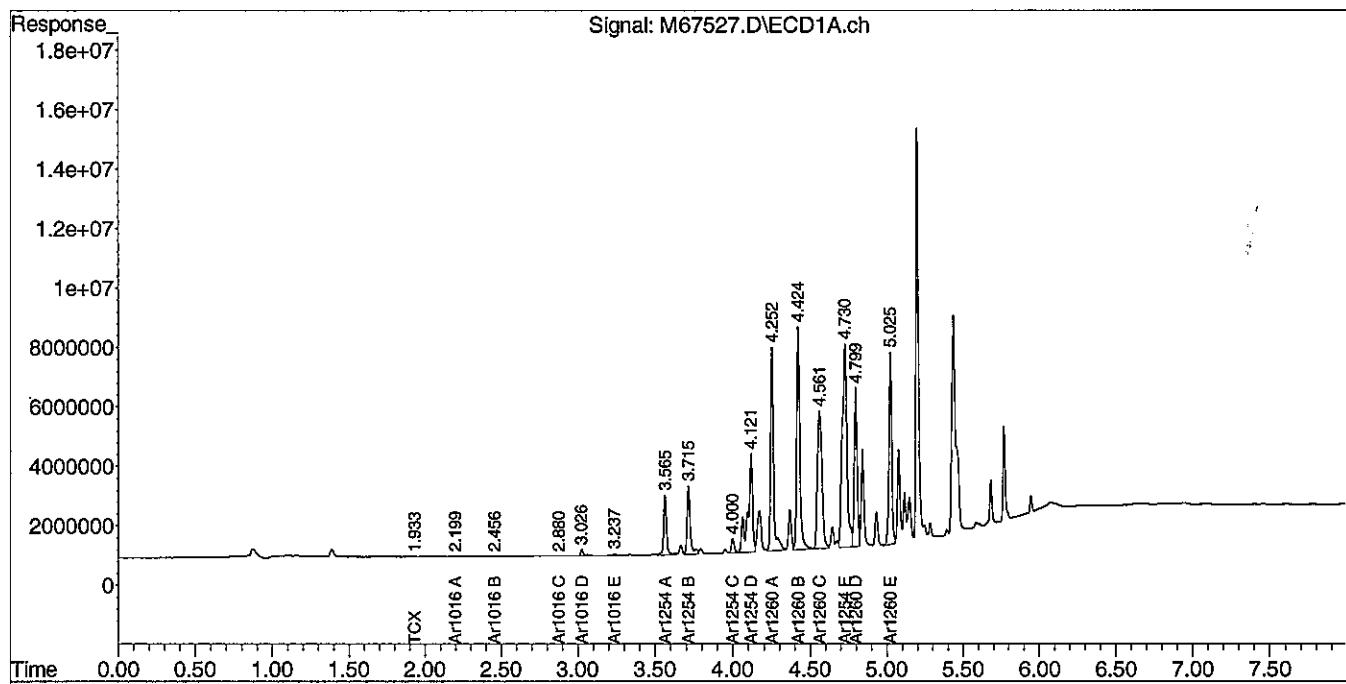
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67527.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 5:21 pm
 Operator : JK
 Sample : 74875-4,1:20000,,A/C
 Misc : SOIL
 ALS Vial : 8 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 17:32:03 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-226 (4-5')

Lab Sample ID: 74875-5
Matrix: Solid
Percent Solid: 84
Dilution Factor: 5900
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	195000	U
PCB-1221	195000	U
PCB-1232	195000	U
PCB-1242	195000	U
PCB-1248	195000	U
PCB-1254	195000	U
PCB-1260	195000	3650000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-5,1:5000,,A/C

Column ID: 0.25 mm

Data File: M67528.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5902.7

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3645900	3288138	10.3	

Column to be used to flag RPD values greater than QC limit of 40%

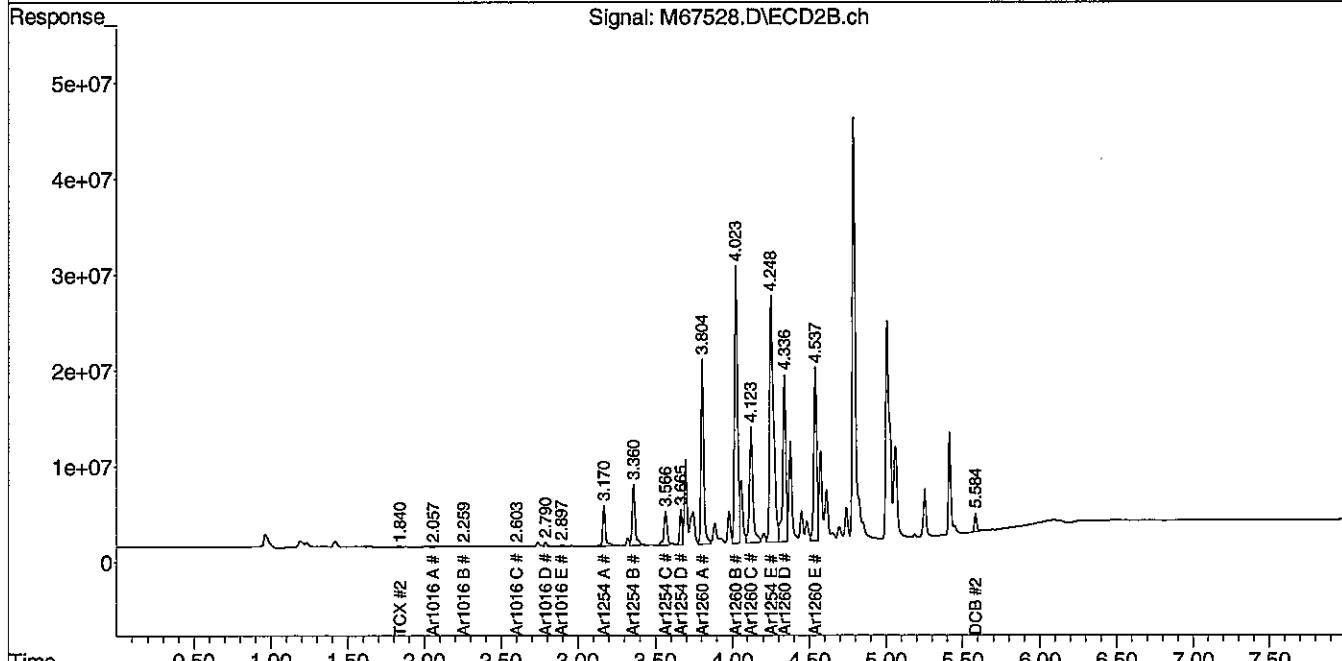
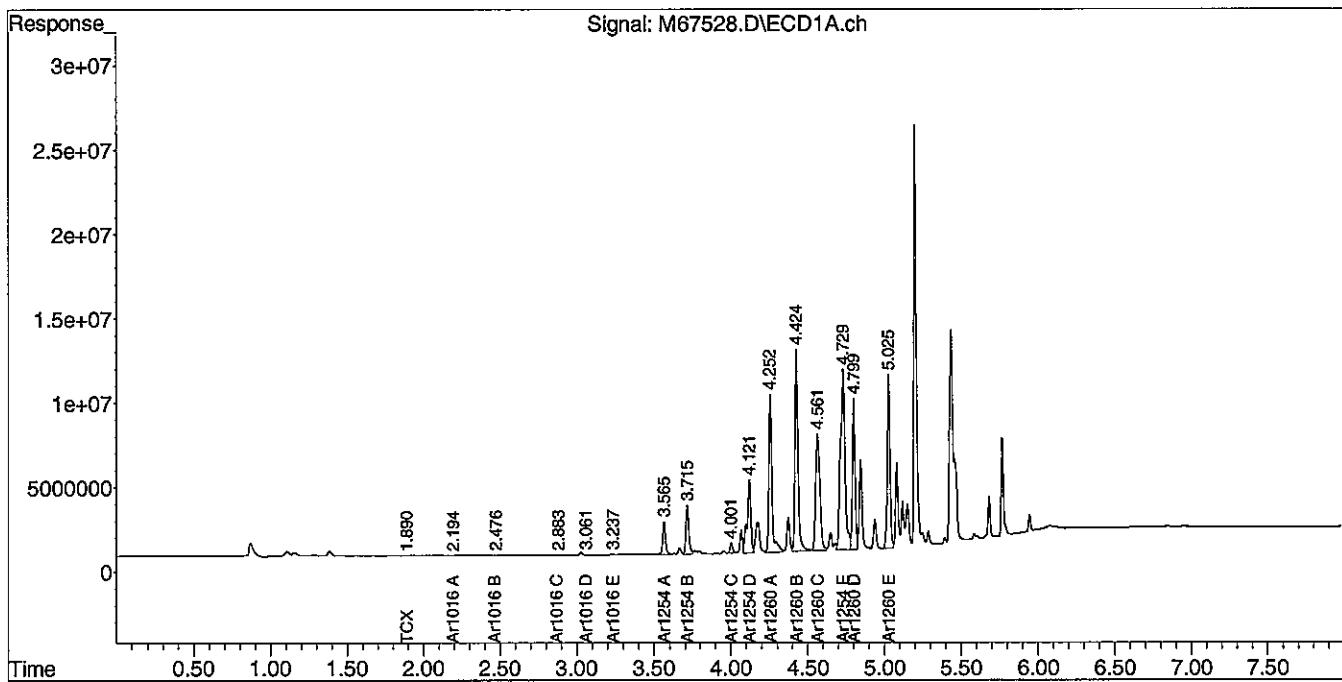
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67528.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 5:31 pm
 Operator : JK
 Sample : 74875-5,1:5000,,A/C
 Misc : SOIL
 ALS Vial : 9 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 17:40:39 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-232 (0-2')

Lab Sample ID: 74875-6
 Matrix: Solid
 Percent Solid: 87
 Dilution Factor: 108
 Collection Date: 02/20/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/22/13
 Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	3560	U
PCB-1221	3560	U
PCB-1232	3560	U
PCB-1242	3560	U
PCB-1248	3560	U
PCB-1254	3560	U
PCB-1260	3560	55300

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-6,1:100,,A/C

Column ID: 0.25 mm

Data File: M67510.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 108.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	55298	50450	9.2	

Column to be used to flag RPD values greater than QC limit of 40%

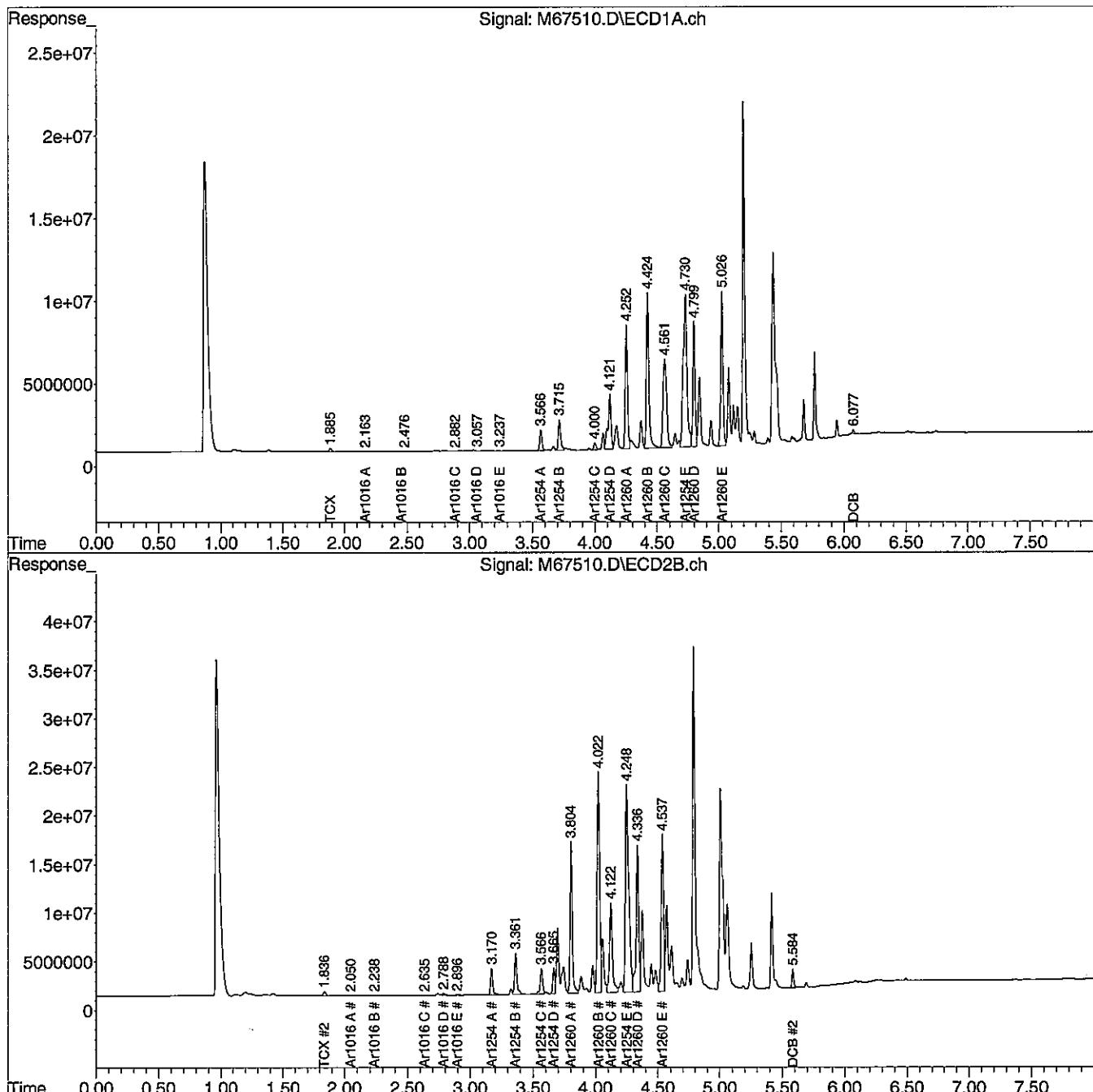
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67510.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 2:29 pm
 Operator : JK
 Sample : 74875-6,1:100,,A/C
 Misc : SOIL
 ALS Vial : 13 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:33 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-232A (2-4')

Lab Sample ID: 74875-7
Matrix: Solid
Percent Solid: 86
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	4490

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	123	%
Decachlorobiphenyl	102	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-7,1:5,,A/C

Column ID: 0.25 mm

Data File: M67511.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.8

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	4486	3825	15.9	

Column to be used to flag RPD values greater than QC limit of 40%

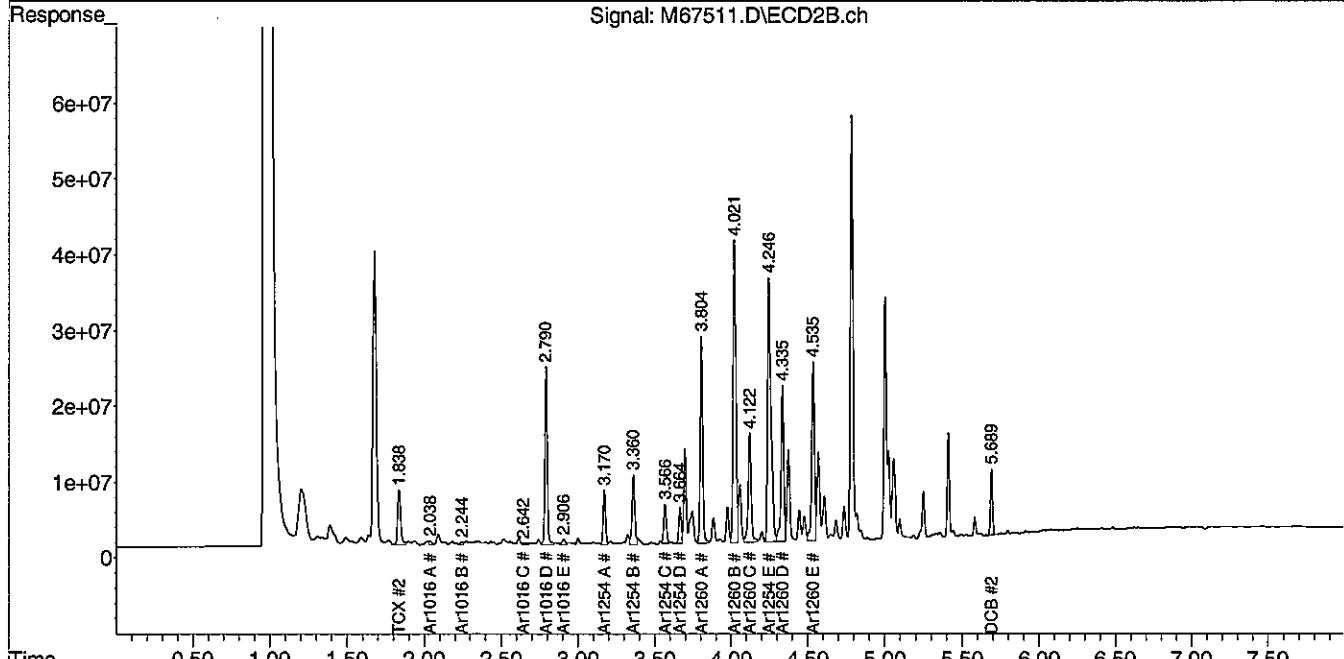
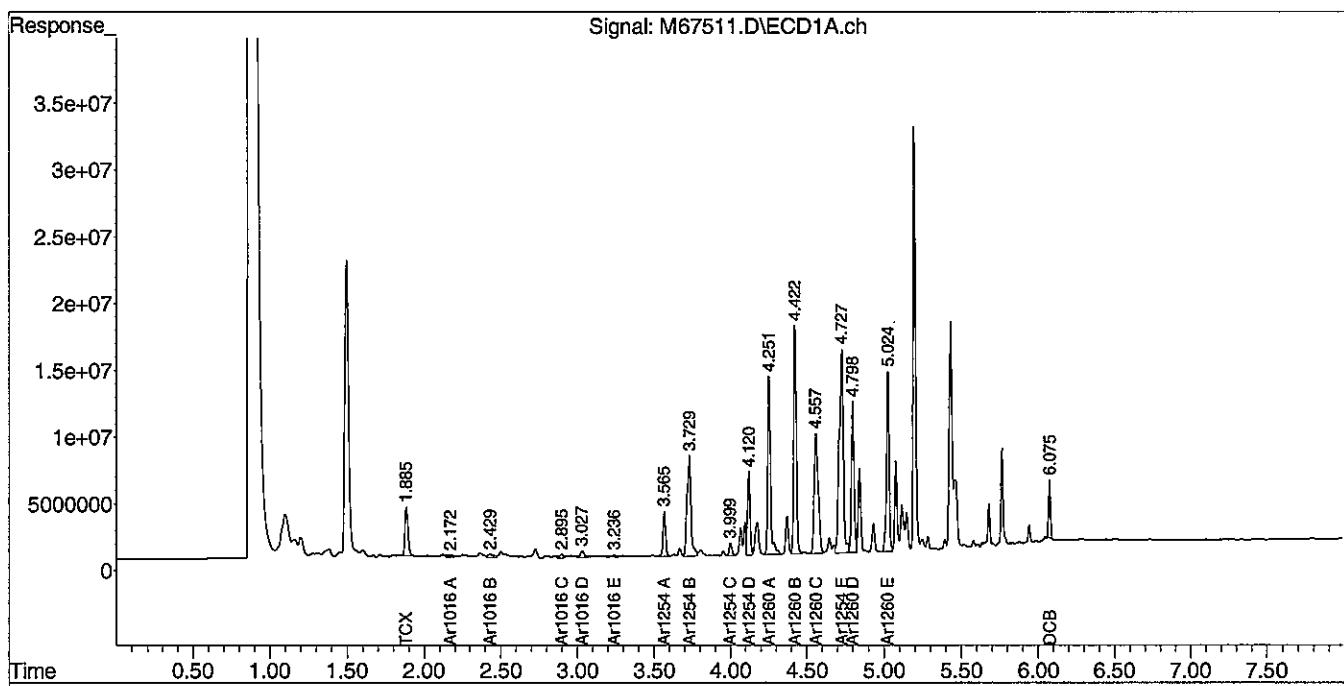
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67511.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 2:39 pm
 Operator : JK
 Sample : 74875-7,1:5,,A/C
 Misc : SOIL
 ALS Vial : 14 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 16:24:35 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-232A (4-6')

Lab Sample ID: 74875-8
Matrix: Solid
Percent Solid: 81
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS			
COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$	
PCB-1016	198		U
PCB-1221	198		U
PCB-1232	198		U
PCB-1242	198		U
PCB-1248	198		U
PCB-1254	198		U
PCB-1260	198		4340

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	98	%	
Decachlorobiphenyl	62	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-8,1:5,,A/C

Column ID: 0.25 mm

Data File: M67514.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 6.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	4343	3695	16.1	

Column to be used to flag RPD values greater than QC limit of 40%

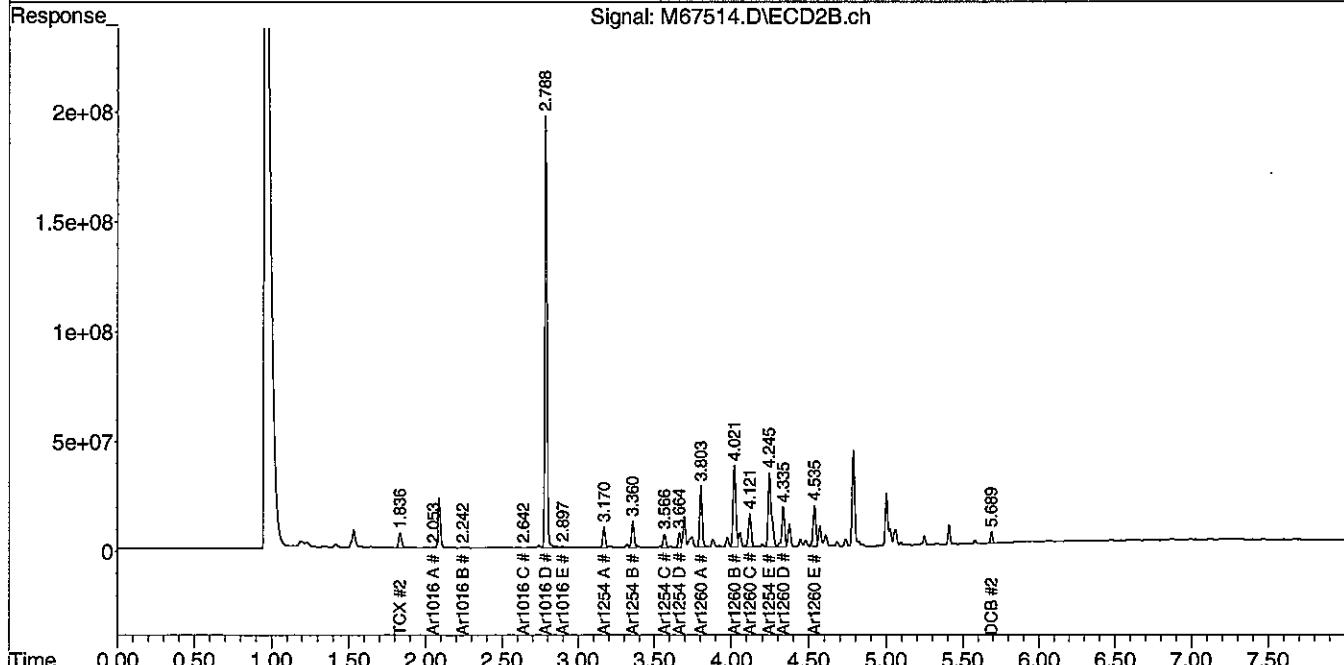
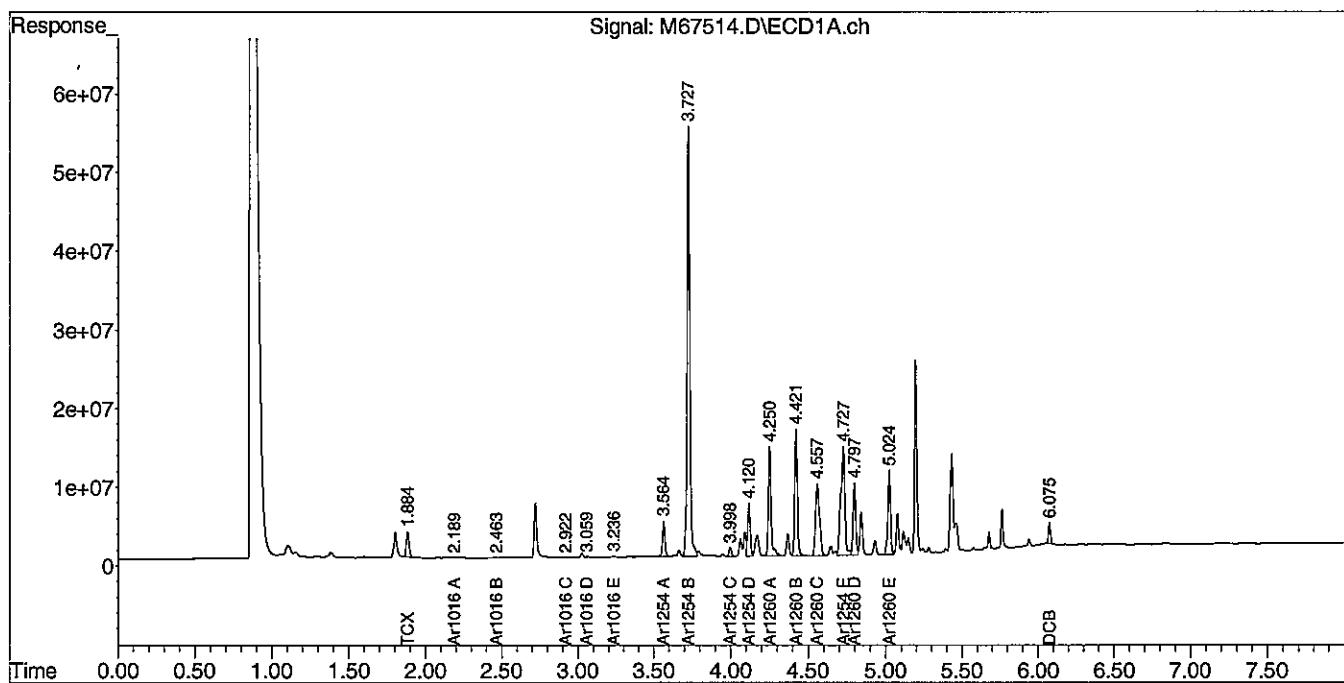
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67514.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 3:09 pm
 Operator : JK
 Sample : 74875-8,1:5,,A/C
 Misc : SOIL
 ALS Vial : 17 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:41 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-232A (6-8')

Lab Sample ID: 74875-9
Matrix: Solid
Percent Solid: 78
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	5090

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	99	%
Decachlorobiphenyl	67	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-9,1:5,,A/C

Column ID: 0.25 mm

Data File: M67515.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 6.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	5091	4104	21.5	

Column to be used to flag RPD values greater than QC limit of 40%

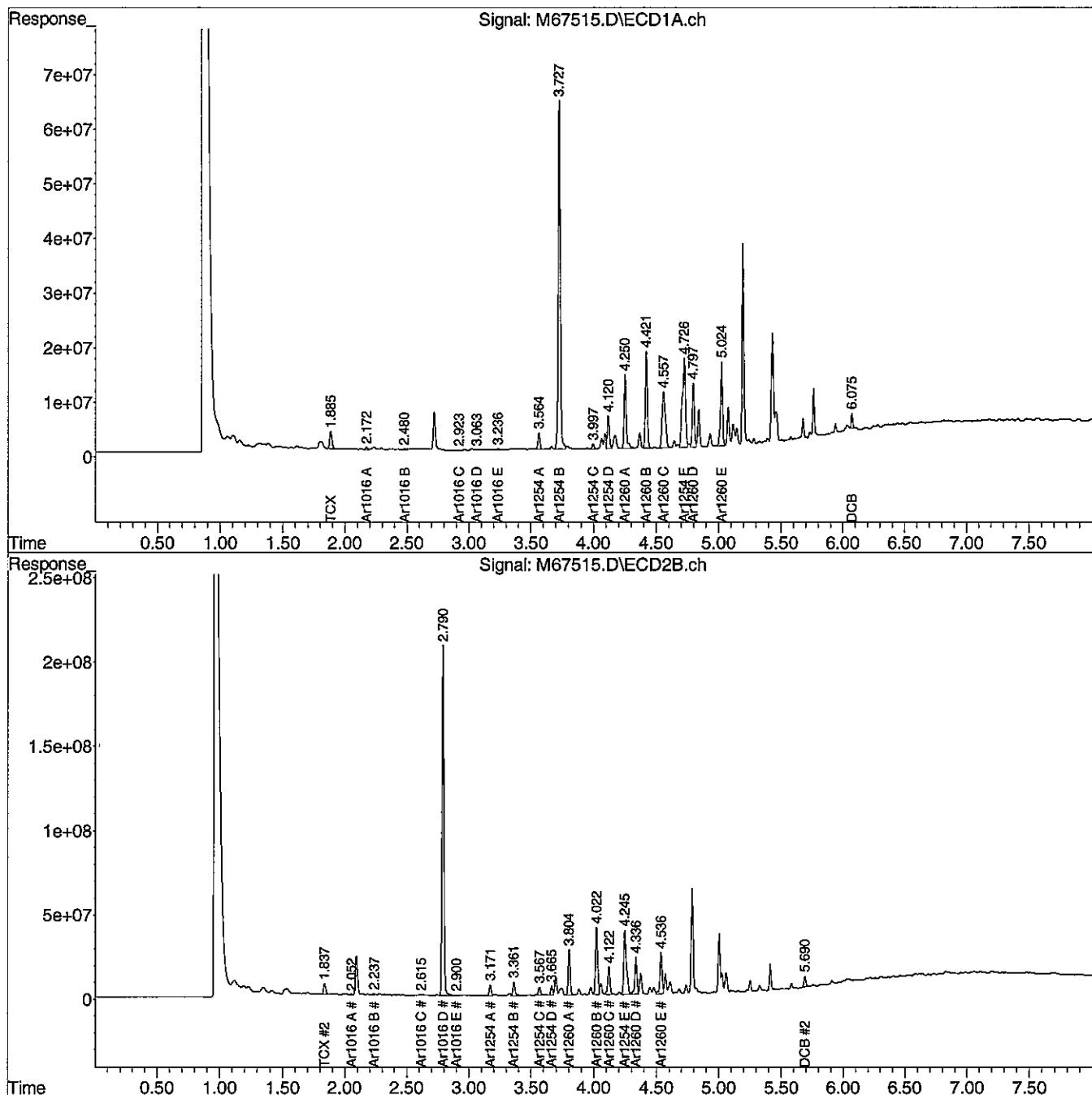
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67515.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 3:19 pm
 Operator : JK
 Sample : 74875-9,1:5,,A/C
 Misc : SOIL
 ALS Vial : 18 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:43 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-232A (8-10')

Lab Sample ID: 74875-10
 Matrix: Solid
 Percent Solid: 89
 Dilution Factor: 1.1
 Collection Date: 02/20/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/22/13
 Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	36	U
PCB-1221	36	U
PCB-1232	36	U
PCB-1242	36	U
PCB-1248	36	U
PCB-1254	36	U
PCB-1260	36	241

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	101	%
Decachlorobiphenyl	76	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-10,,A/C

Column ID: 0.25 mm

Data File: M67516.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	241	202	17.8	

Column to be used to flag RPD values greater than QC limit of 40%

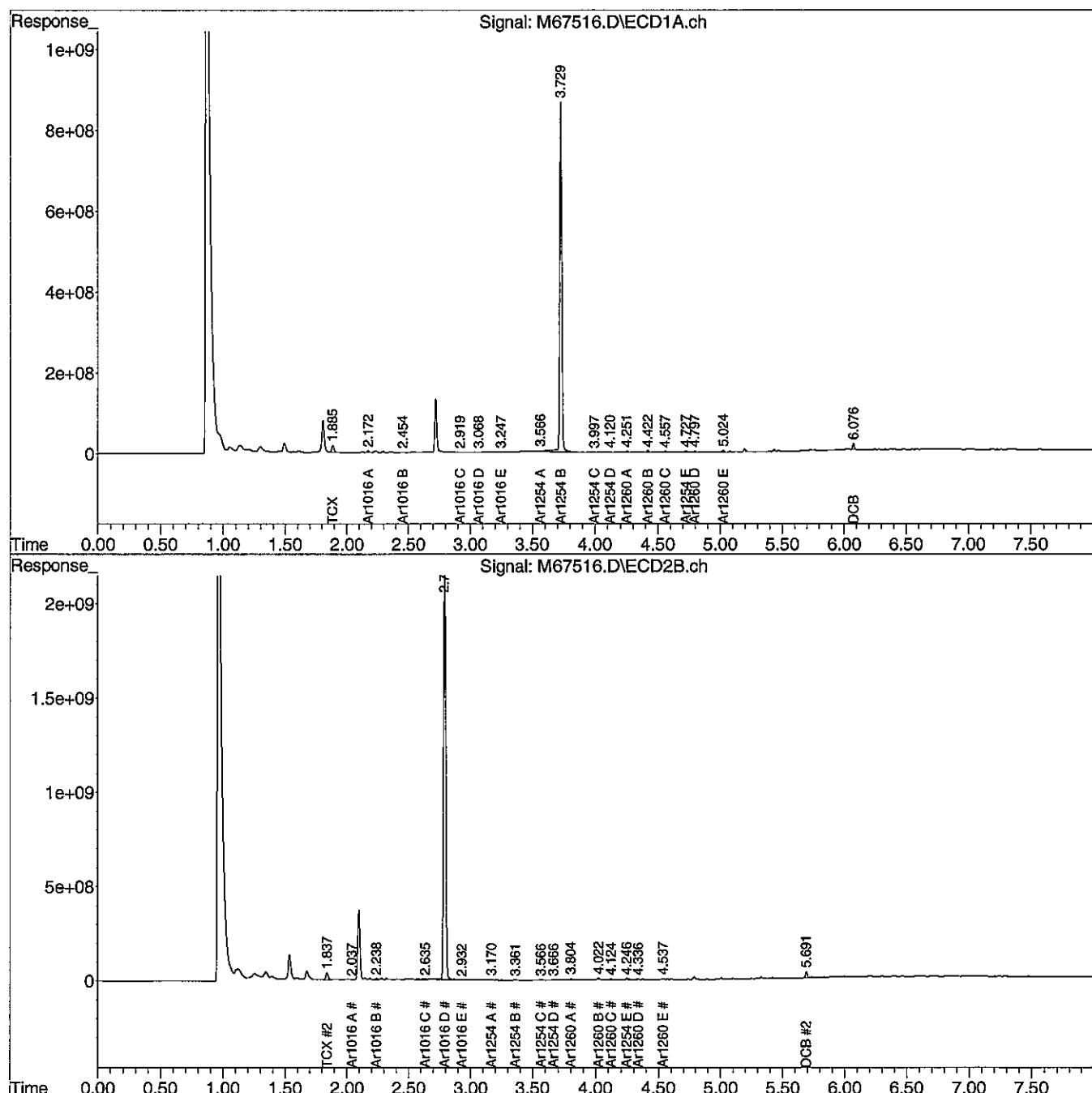
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67516.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 3:30 pm
 Operator : JK
 Sample : 74875-10,, A/C
 Misc : SOIL
 ALS Vial : 19 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:45 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-233 (0-2')

Lab Sample ID: 74875-11
Matrix: Solid
Percent Solid: 80
Dilution Factor: 12
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	396	U
PCB-1221	396	U
PCB-1232	396	U
PCB-1242	396	U
PCB-1248	396	U
PCB-1254	396	U
PCB-1260	396	6300

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	108	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-11,1:10,,A/C

Column ID: 0.25 mm

Data File: M67517.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 12.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	6304	5298	17.3	

Column to be used to flag RPD values greater than QC limit of 40%

* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67517.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 3:40 pm
 Operator : JK
 Sample : 74875-11,1:10,,A/C
 Misc : SOIL
 ALS Vial : 20 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e

Quant Time: Feb 27 16:32:50 2013

Quant Method : C:\msdchem\1\METHODS\PCB020513.M

Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254

QLast Update : Wed Feb 27 09:33:47 2013

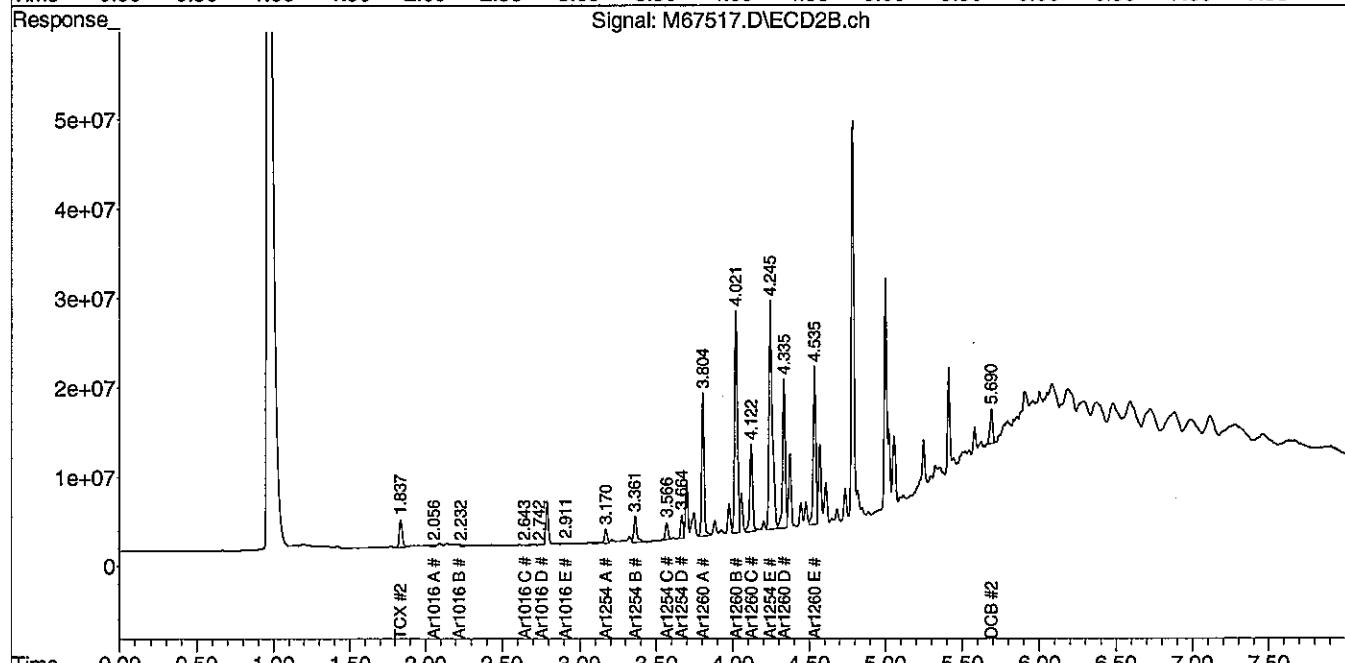
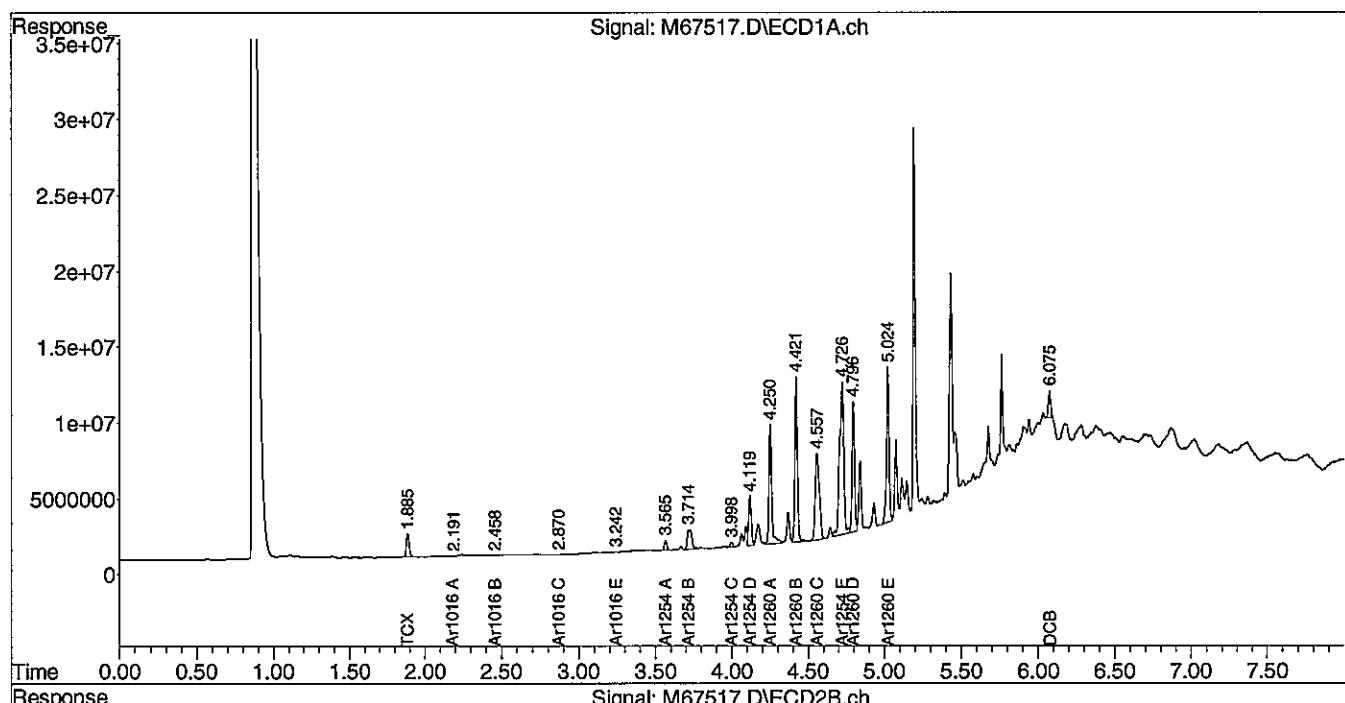
Response via : Initial Calibration

Integrator: ChemStation

Volume Inj. : 2 uL

Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides

Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-233 (2-4')

Lab Sample ID: 74875-12
 Matrix: Solid
 Percent Solid: 90
 Dilution Factor: 1.0
 Collection Date: 02/20/13
 Lab Receipt Date: 02/20/13
 Extraction Date: 02/22/13
 Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	1020

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	96	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-12,,A/C

Column ID: 0.25 mm

Data File: M67518.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1022	865	16.6	

Column to be used to flag RPD values greater than QC limit of 40%

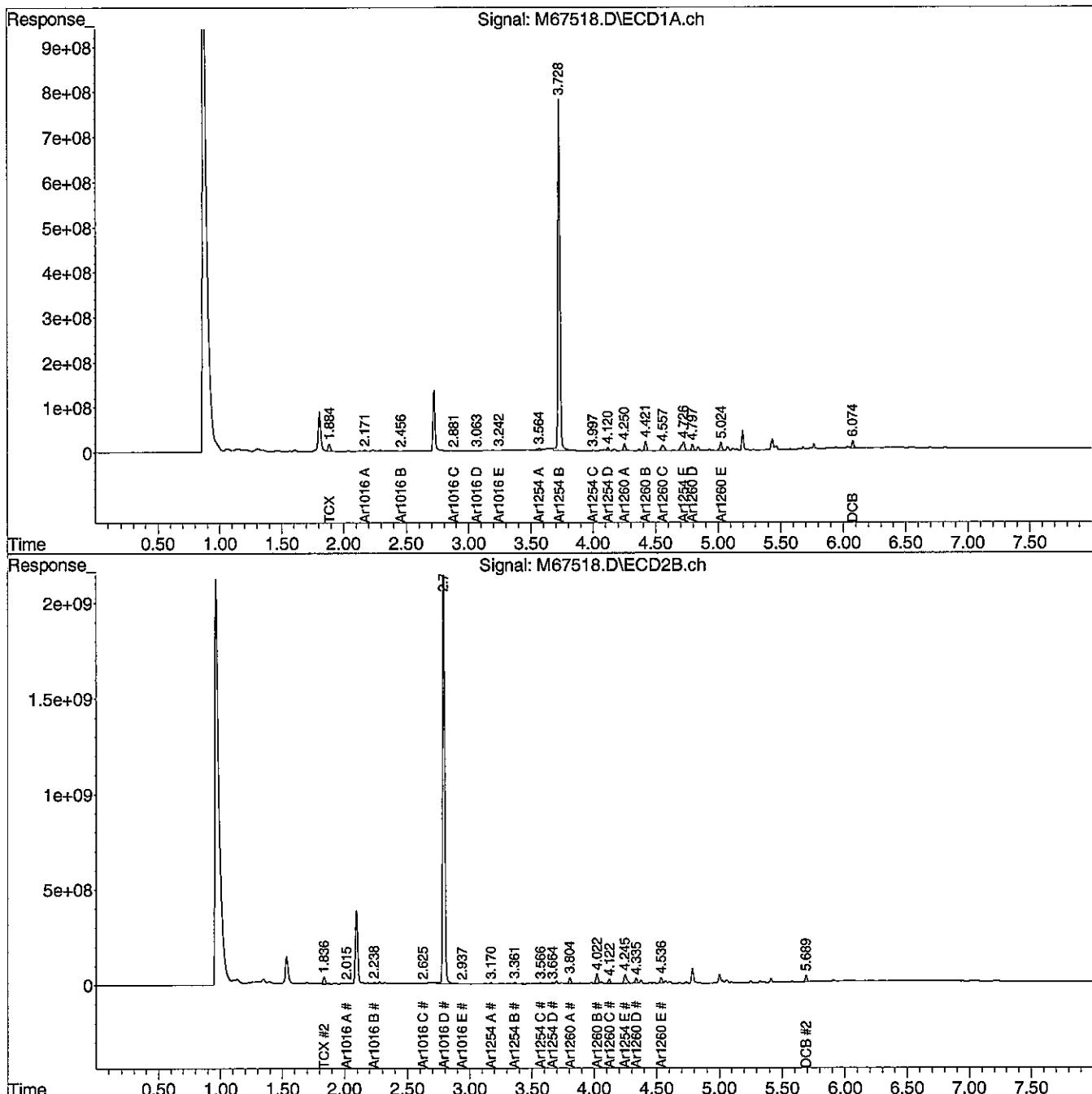
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
Data File : M67518.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 27 Feb 2013 3:50 pm
Operator : JK
Sample : 74875-12,,A/C
Misc : SOIL
ALS Vial : 21 Sample Multiplier: 1

```
Integration File signal 1: events.e
Integration File signal 2: events2.e
Quant Time: Feb 27 16:02:02 2013
Quant Method : C:\msdchem\1\METHODS\PCB020513.M
Quant Title  : SW-846 METHOD 8082 Aroclor 1016/1260/1254
QLast Update : Wed Feb 27 09:33:47 2013
Response via : Initial Calibration
Integrator: ChemStation
```

Volume Inj. : 2 uL
Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-233 (4-6')

Lab Sample ID: 74875-13
Matrix: Solid
Percent Solid: 91
Dilution Factor: 1.0
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	267

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	106	%
Decachlorobiphenyl	78	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-13,,A/C

Column ID: 0.25 mm

Data File: M67519.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	267	229	15.3	

Column to be used to flag RPD values greater than QC limit of 40%

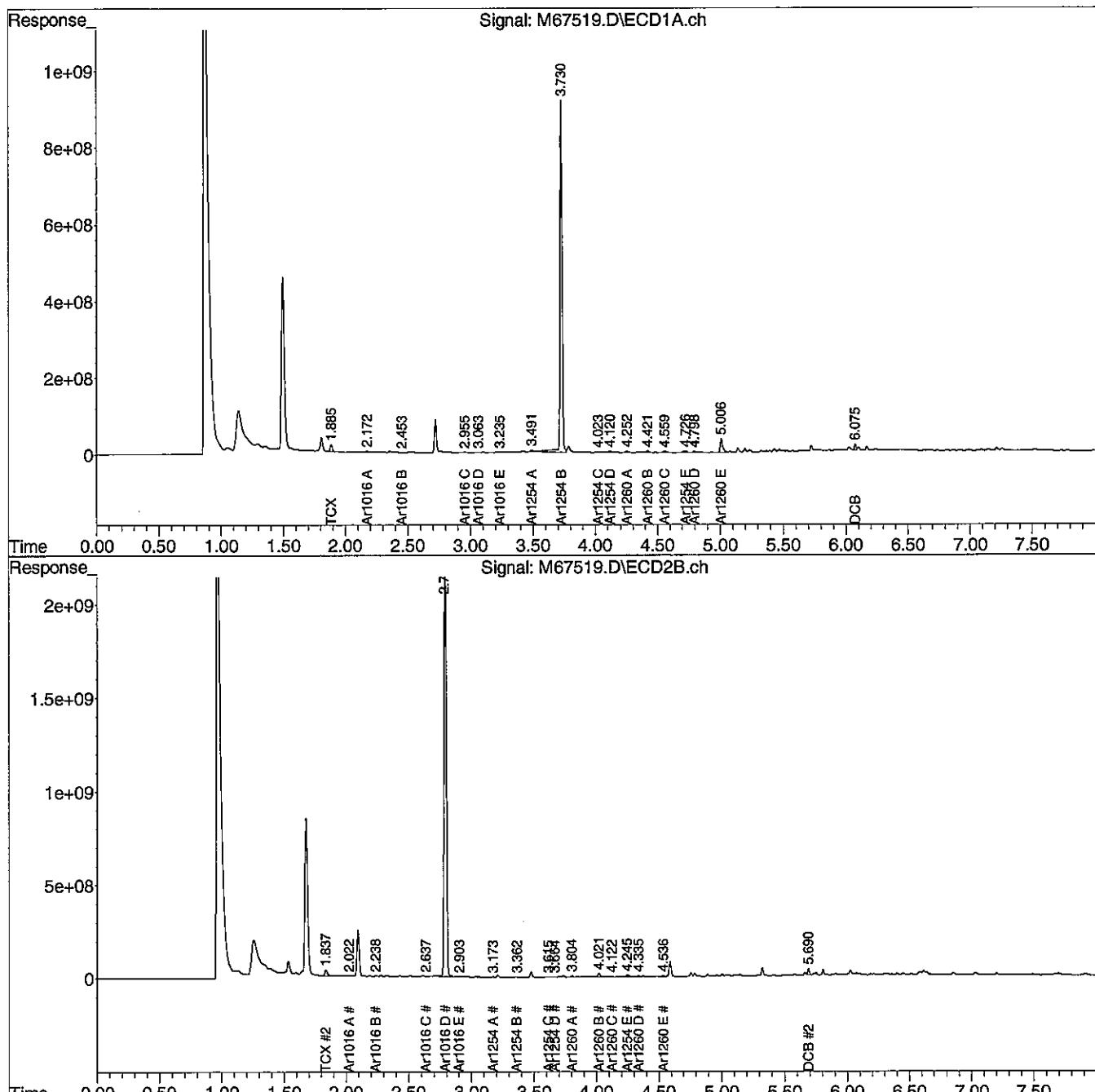
* Values outside QC limits

Comments: _____

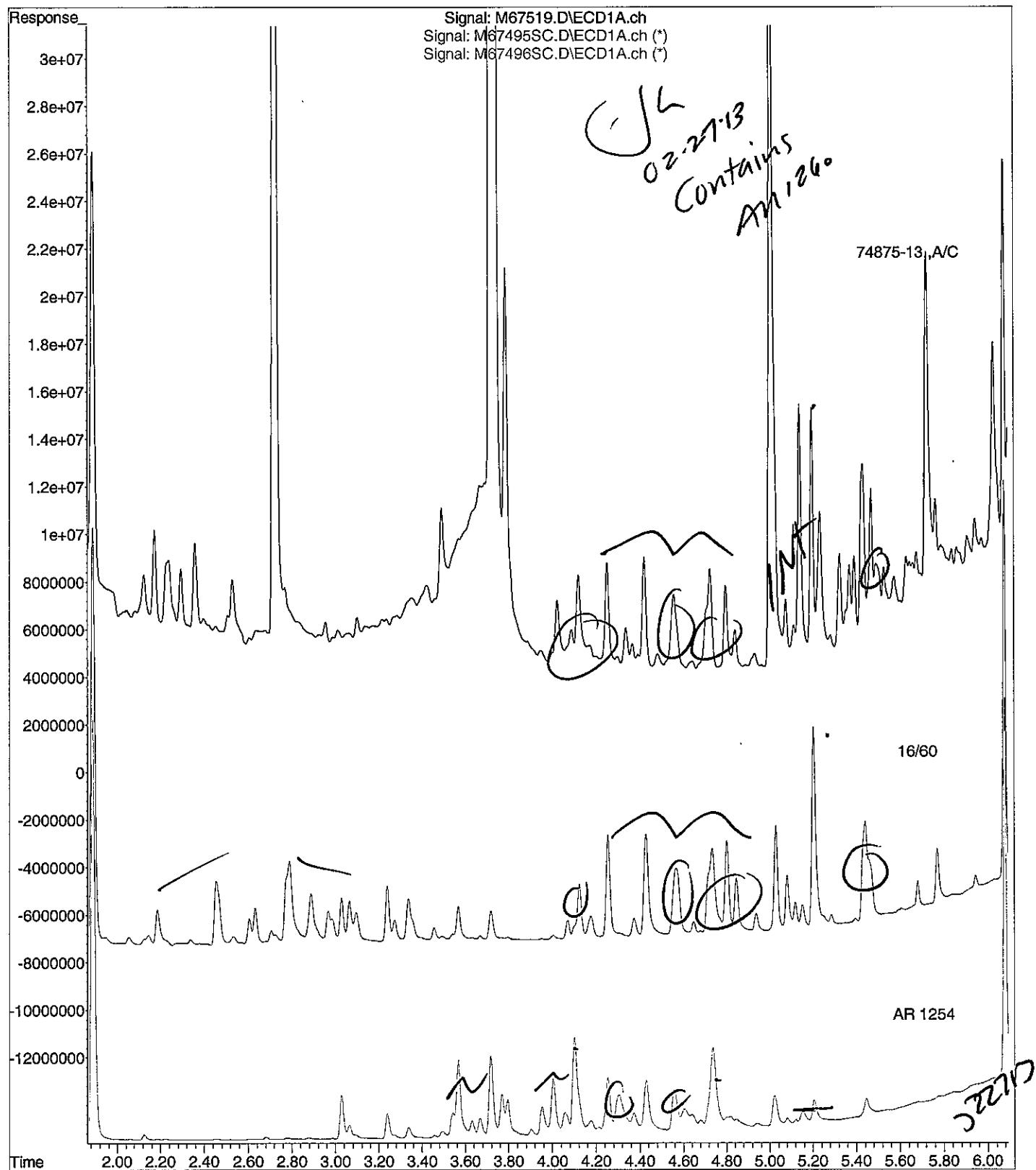
Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67519.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 4:00 pm
 Operator : JK
 Sample : 74875-13,,A/C
 Misc : SOIL
 ALS Vial : 22 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 16:09:38 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022713-M\M67519.D
Operator : JK
Acquired : 27 Feb 2013 4:00 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74875-13,,A/C
Misc Info : SOIL
Vial Number: 22



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-233 (6-8')

Lab Sample ID: 74875-14
Matrix: Solid
Percent Solid: 89
Dilution Factor: 1.0
Collection Date: 02/20/13
Lab Receipt Date: 02/20/13
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	404

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	93	%
Decachlorobiphenyl	88	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG:

GC Column #1: STX-CLPesticides I

Sample: 74875-14,,A/C

Column ID: 0.25 mm

Data File: M67520.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	404	364	10.3	

Column to be used to flag RPD values greater than QC limit of 40%

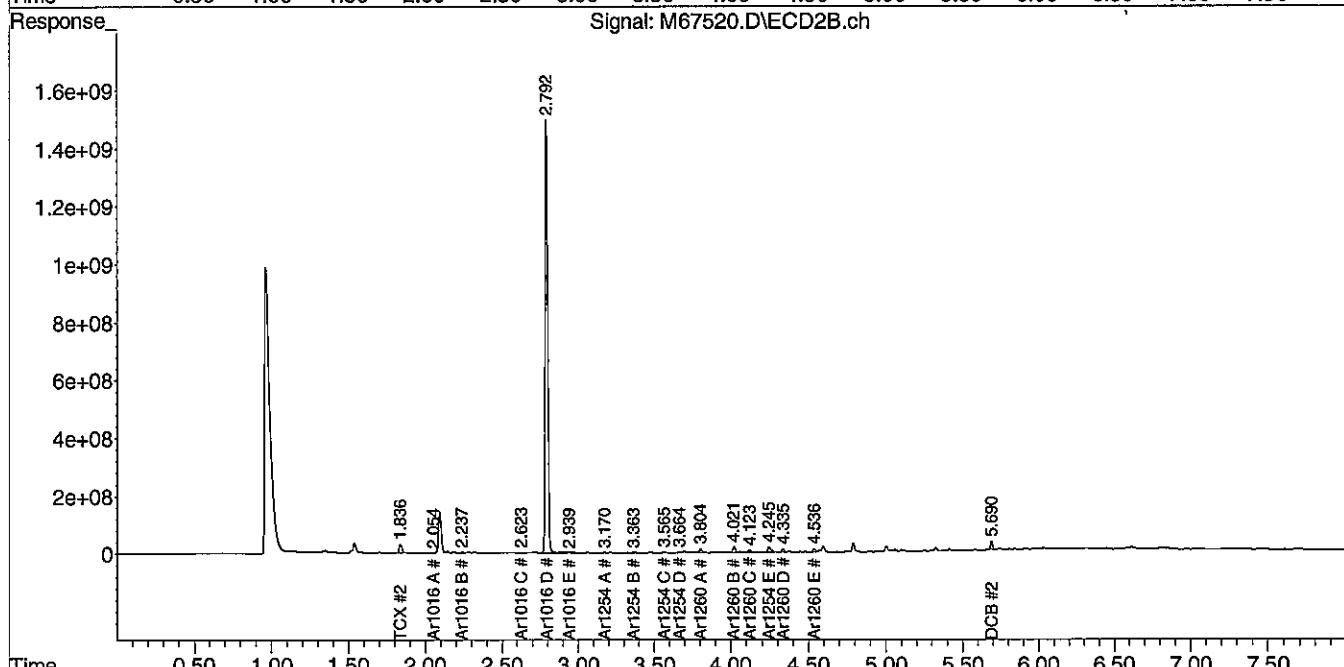
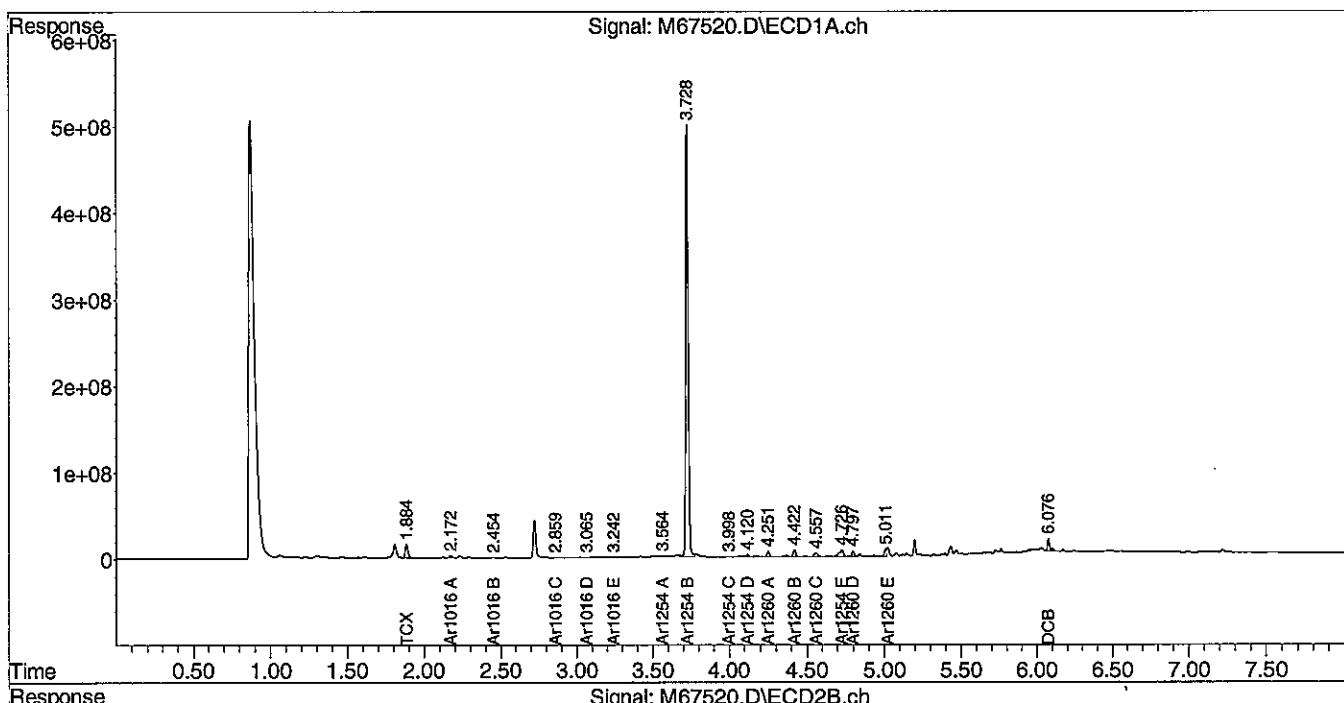
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67520.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 4:10 pm
 Operator : JK
 Sample : 74875-14,,A/C
 Misc : SOIL
 ALS Vial : 23 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 16:23:15 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-234 (0-2")

Lab Sample ID:	74875-15
Matrix:	Solid
Percent Solid:	88
Dilution Factor:	52
Collection Date:	02/20/13
Lab Receipt Date:	02/20/13
Extraction Date:	02/22/13
Analysis Date:	02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	1720	U
PCB-1221	1720	U
PCB-1232	1720	U
PCB-1242	1720	U
PCB-1248	1720	U
PCB-1254	1720	U
PCB-1260	1720	25100

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74875

GC Column #1: STX-CLPesticides I

Sample: 74875-15,1:50,,A/C

Column ID: 0.25 mm

Data File: M67521.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 52.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	25113	22324	11.8	

Column to be used to flag RPD values greater than QC limit of 40%

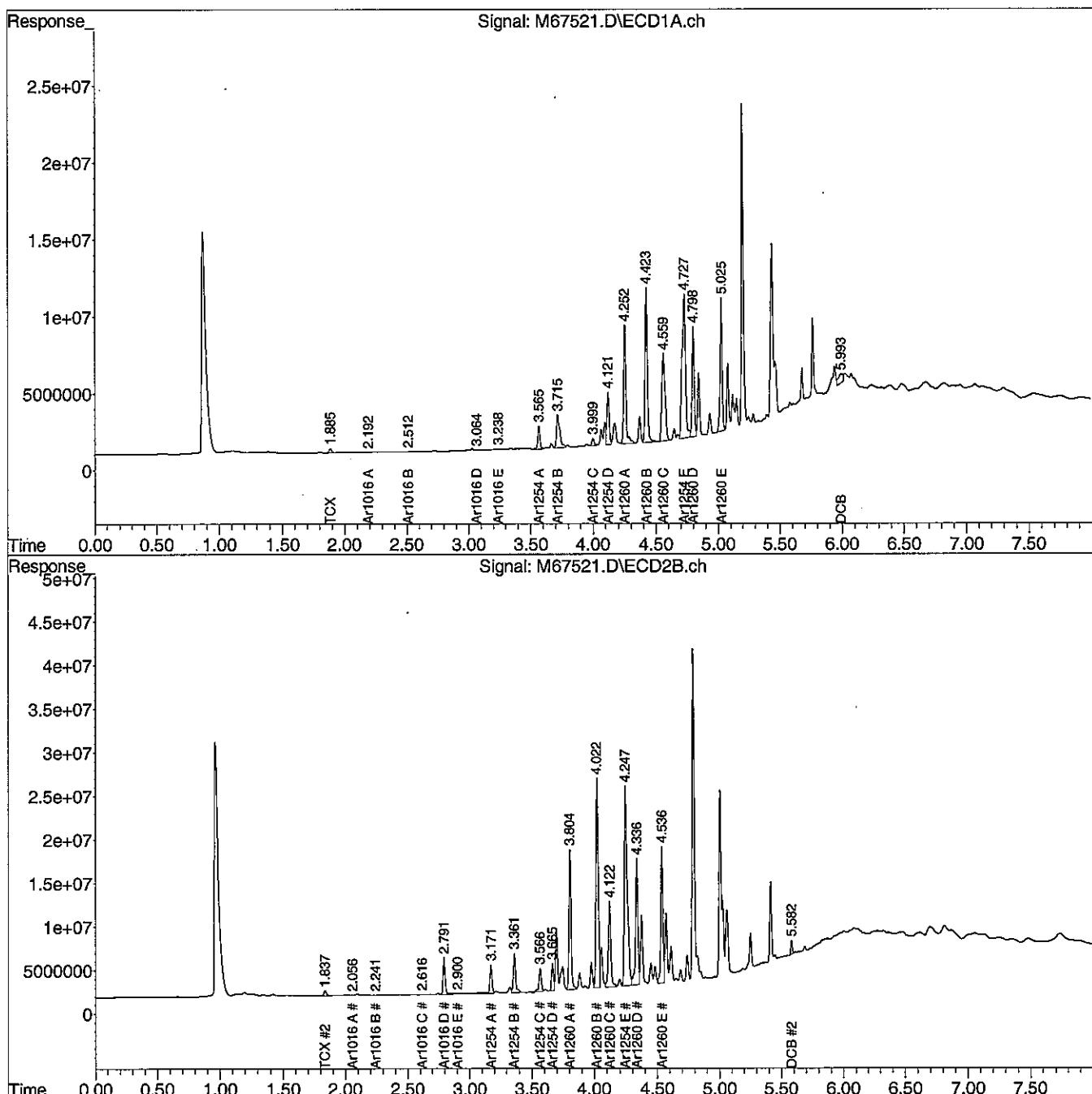
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67521.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 4:20 pm
 Operator : JK
 Sample : 74875-15,1:50,,A/C
 Misc : SOIL
 ALS Vial : 24 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 16:33:01 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



PCB
QC FORMS

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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022213PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	69	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022613-M\
 Data File : M67454B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:46 am *Jn 022713*
 Operator : JK
 Sample : B022213PSOX,, A/C (Sig #1); B022113PSOX,, A/C (Sig #2)
 Misc : SOIL
 ALS Vial : 38 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e

Quant Time: Feb 27 12:47:07 2013

Quant Method : C:\msdchem\1\METHODS\PCB020513.M

Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254

QLast Update : Wed Feb 27 09:33:47 2013

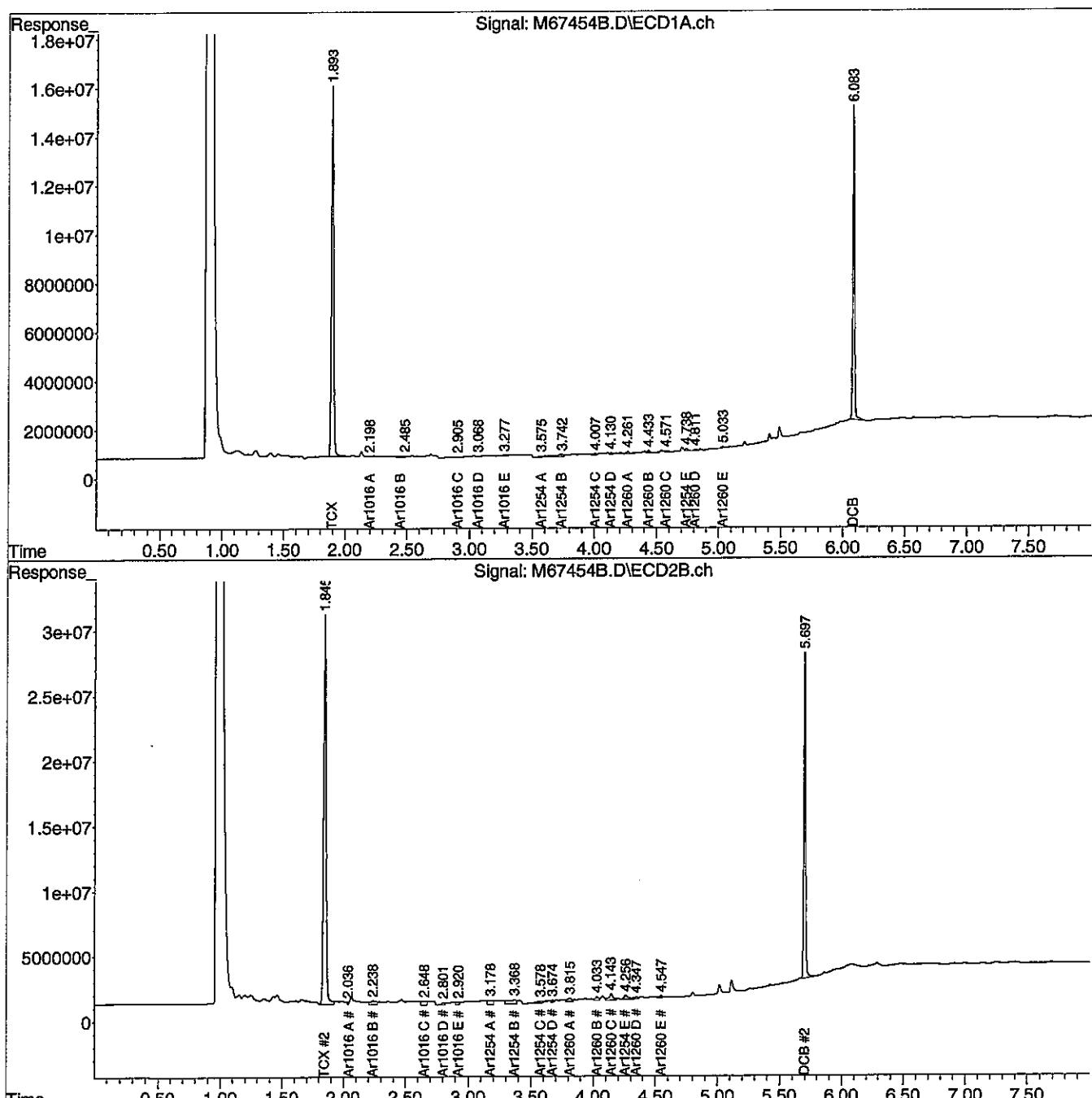
Response via : Initial Calibration

Integrator: ChemStation

Volume Inj. : 2 uL

Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides

Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022213PSOX RR
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	75	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

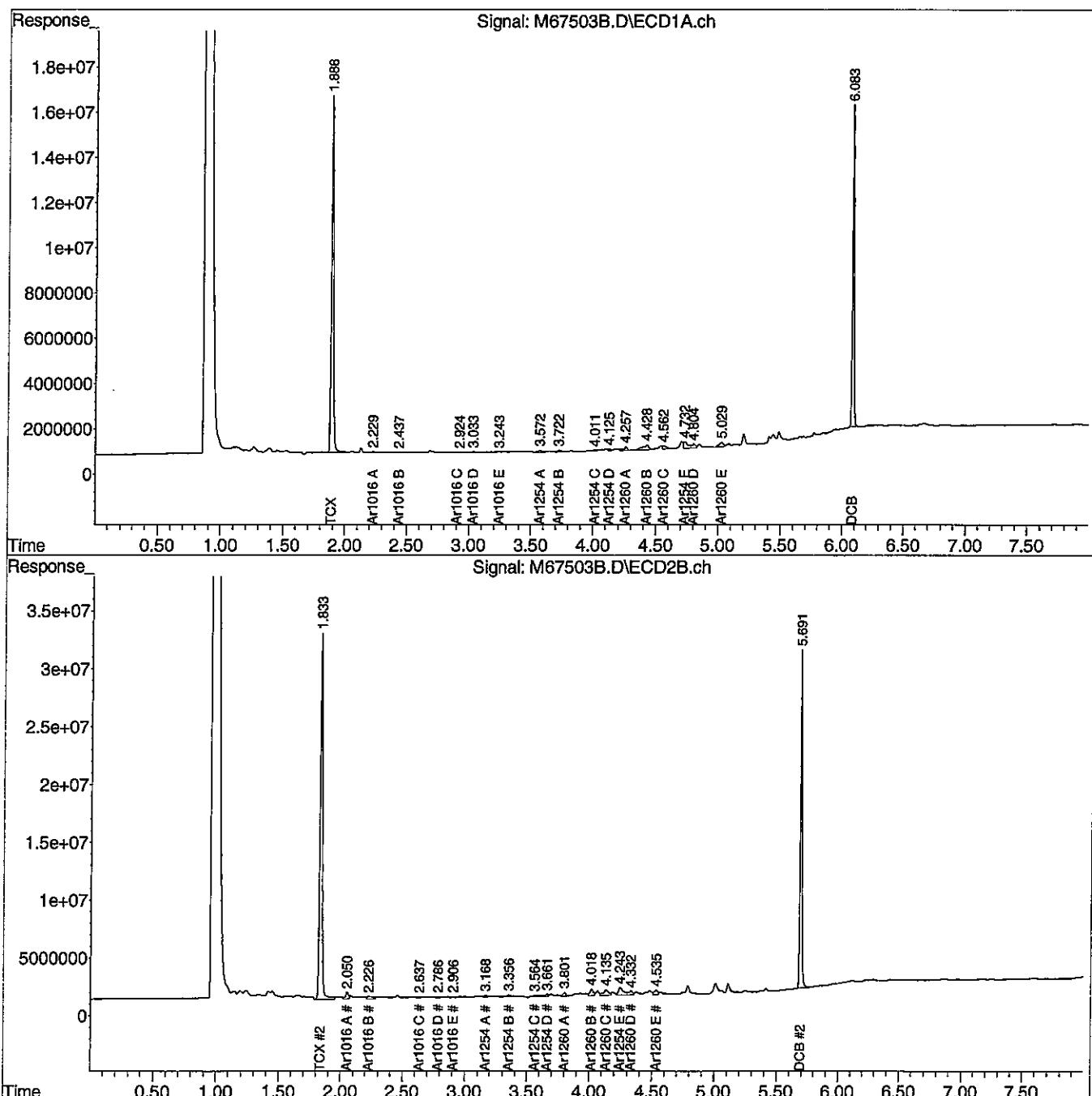
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67503B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 1:19 pm
 Operator : JK
 Sample : B022213PSOX,RR2,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 15:42:19 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

February 27, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022213PSOX RR
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/27/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	71	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

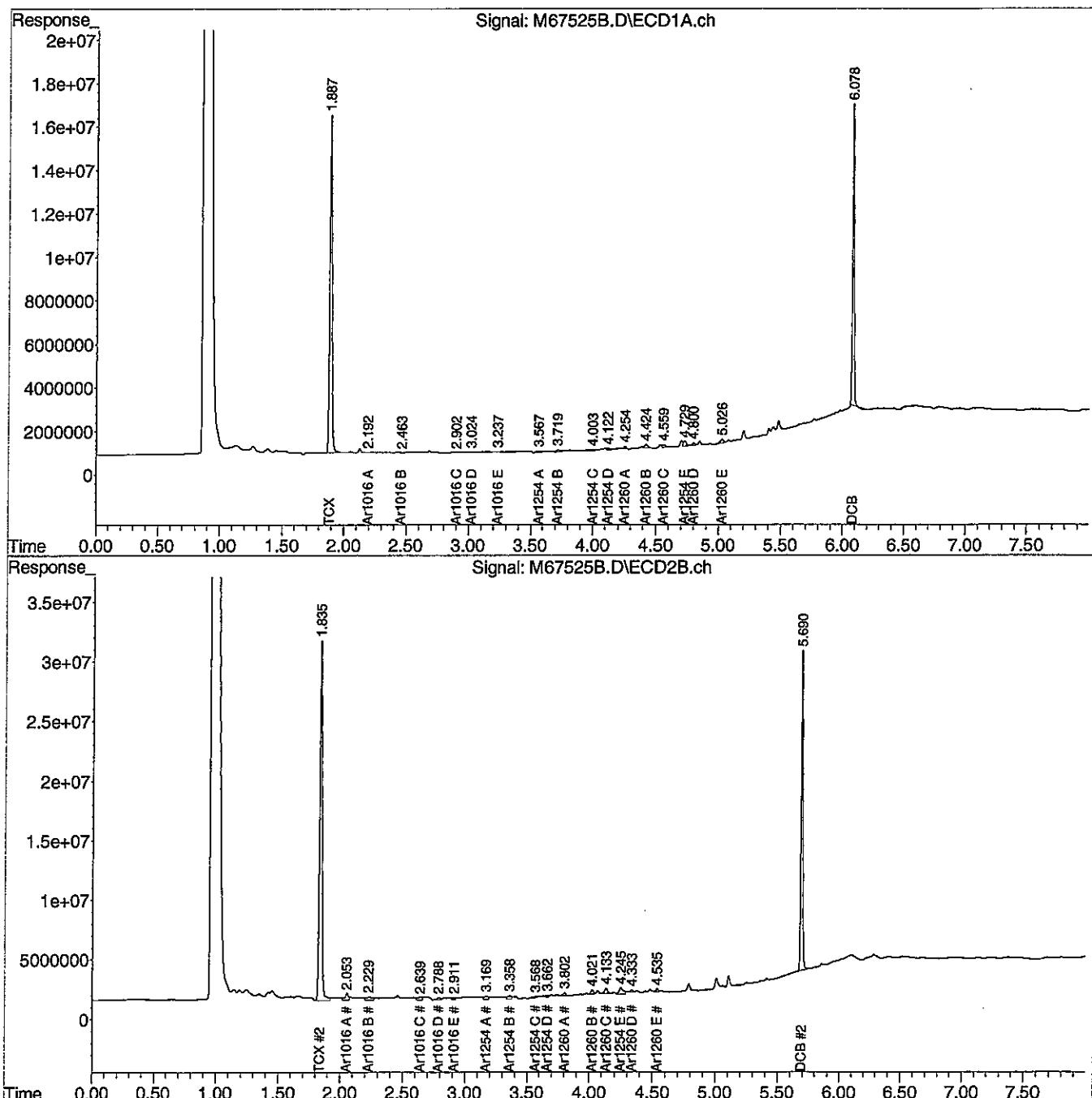
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022713-M\
 Data File : M67525B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 27 Feb 2013 5:01 pm
 Operator : JK
 Sample : B022213PSOX,RR3,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 27 17:15:15 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Feb 27 09:33:47 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74875

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74875

Column ID: 0,25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

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SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74875

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022213PSOX,,A/C

Spike: L022213PSOX,,A/C

Spike duplicate: LD022213PSOX,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	178	89		175	87		2.0	
PCB 1260	200	200	60	130	30	0	179	90		173	87		3.3	
PCB 1016 #2	200	200	65	140	30	0	173	86		166	83		4.0	
PCB 1260 #2	200	200	60	130	30	0	172	86		166	83		3.7	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: 74875-7,1:5,,A/C

Spike: 74875-7,MS,1:5,,A/C

Spike duplicate: 74875-7,MSD,1:5,,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP		RPD	#
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#		
PCB 1016	229	224	65	140	30	0	D	D		D	D			
PCB 1260	229	224	60	130	30	4486	5470	429	*	3332	-514	*	48.6	*
PCB 1016 #2	229	224	65	140	30	0	D	D		D	D			
PCB 1260 #2	229	224	60	130	30	3825	4650	360	*	2862	-430	*	47.6	*

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form



193 Commerce way, Suite E
Portsmouth, NH 03801
(800) 929-9906
(603) 430-2151 Fax
(603) 438-5111

		Project Name: MAINE ENERGY Project #: 12-3359.1		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax																																																																																									
<p style="text-align: center;">For Analytics Use Only</p> <p>Samples were:</p> <ol style="list-style-type: none"> 1) Shipped or hand-delivered <input checked="" type="checkbox"/> 2) Temperature (°C): <u>0 - 30.70</u> 3) Received in good condition: <input checked="" type="checkbox"/> N/A 4) pH checked by: <u>N/A</u> 5) Labels checked by: <u>N/A</u> 																																																																																															
<p>Matrix Key:</p> <ul style="list-style-type: none"> C = Concrete GW = Groundwater WP = Wrap SW = Surface Water HNO3 = HNO3 Oil = Oil Other = Other Extract = Extract 																																																																																															
<table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="text-align: left; width: 15%;">Sample Identification</th> <th style="text-align: left; width: 15%;">Sample Date</th> <th style="text-align: left; width: 15%;">Sample Time</th> <th style="text-align: left; width: 15%;">Field Filtered Y or N</th> <th style="text-align: left; width: 15%;">Preservative Code:</th> <th style="text-align: left; width: 15%;">Please <input checked="" type="checkbox"/> in preservation code here</th> <th style="text-align: left; width: 15%;">Full or Ranges only TETPH</th> <th style="text-align: left; width: 15%;">Full or Ranges only RCRA8 PPI3 TAL23 Other**</th> </tr> </thead> <tbody> <tr> <td>SE-SS-225 (4-6)</td> <td>2-19</td> <td>1741</td> <td>N</td> <td>X X</td> <td>S</td> <td>1</td> <td>74875-1</td> </tr> <tr> <td>SE-SS-225 (6-7.5)</td> <td>2-19</td> <td>1743</td> <td>N</td> <td>X X</td> <td>S</td> <td>1</td> <td>-2</td> </tr> <tr> <td>SE-SS-226 (8-3)</td> <td>2-20</td> <td>835</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-3</td> </tr> <tr> <td>SE-SS-226 (3-4)</td> <td>2-20</td> <td>835</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-4</td> </tr> <tr> <td>SE-SS-224 (4-5)</td> <td>2-20</td> <td>910</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-5</td> </tr> <tr> <td>SE-SS-232 (5-6)</td> <td>2-20</td> <td>950</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-6</td> </tr> <tr> <td>SE-SS-232A (3-4)</td> <td>2-20</td> <td>1020</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-7</td> </tr> <tr> <td>SE-SS-232A (4-6)</td> <td>2-20</td> <td>1025</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-8</td> </tr> <tr> <td>SE-SS-232A (6-8)</td> <td>2-20</td> <td>1035</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-9</td> </tr> <tr> <td>SE-SS-232A (8-10)</td> <td>2-20</td> <td>1040</td> <td>N</td> <td>X</td> <td>S</td> <td>1</td> <td>-10</td> </tr> </tbody> </table>								Sample Identification	Sample Date	Sample Time	Field Filtered Y or N	Preservative Code:	Please <input checked="" type="checkbox"/> in preservation code here	Full or Ranges only TETPH	Full or Ranges only RCRA8 PPI3 TAL23 Other**	SE-SS-225 (4-6)	2-19	1741	N	X X	S	1	74875-1	SE-SS-225 (6-7.5)	2-19	1743	N	X X	S	1	-2	SE-SS-226 (8-3)	2-20	835	N	X	S	1	-3	SE-SS-226 (3-4)	2-20	835	N	X	S	1	-4	SE-SS-224 (4-5)	2-20	910	N	X	S	1	-5	SE-SS-232 (5-6)	2-20	950	N	X	S	1	-6	SE-SS-232A (3-4)	2-20	1020	N	X	S	1	-7	SE-SS-232A (4-6)	2-20	1025	N	X	S	1	-8	SE-SS-232A (6-8)	2-20	1035	N	X	S	1	-9	SE-SS-232A (8-10)	2-20	1040	N	X	S	1	-10
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<p>Comments, Additional Analyses, or Special Instructions:</p> <p>** List requested metals here *Fee may apply</p> <p>Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.</p>																																																																																															
<p>Email Results to:</p> <hr/>				<p>Report Type: <input type="checkbox"/> MCP* <input type="checkbox"/> Level II* <input type="checkbox"/> State: NH <input type="checkbox"/> State Standard:</p> <p><input type="checkbox"/> CTRCP* <input type="checkbox"/> Level III* <input type="checkbox"/> ME <input type="checkbox"/> (eg. S-1 or GW-1)</p> <p><input type="checkbox"/> DOD* <input type="checkbox"/> Level IV* <input type="checkbox"/> CT <input type="checkbox"/> EDD Required: Y/N <input checked="" type="checkbox"/> RI</p> <p><input checked="" type="checkbox"/> Standard <input type="checkbox"/> Other: _____ <input type="checkbox"/> Type: _____</p>																																																																																											
<p>Turnaround Time (TAT)</p> <p><input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> 4 Days* <input checked="" type="checkbox"/> 5 Days <input type="checkbox"/> Standard (8-10 business days)</p> <p>*Fee may apply; lab approval required</p> <p>Sampler Name (Print): <u>John Gresser</u></p>				<p>Date: <u>2-20</u> Time: <u>131</u> Received By: <u>John Gresser</u></p> <p>Date: <u>2-20</u> Time: <u>131</u> Received By: <u>John Gresser</u></p>																																																																																											
<p>Relinquished By: Sampler: <u>John Gresser</u> Relinquished By: <u>John Gresser</u></p>																																																																																															

Chain Of Custody Form

Analytics		environmental laboratory LLC		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax		For Analytics Use Only	
Project Name: MAINE ENERGY		Preservation Code: Report to: SUMMIT ENERGY		Circle and/or Write Required Analysis Followed by Preservation Code Please circle in preservation code here		Samples were: 1) Shipped or Hand-delivered 2) Temperature (°C): 0.3-0.70 3) Received in good condition: Y or N 4) pH checked by: NA 5) Labels checked by: 12/20/13			
Project #: 12-32591		Preservation Key: A = HCl B = 4°C C = Unpres D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other		Metals: RCRA PPL13 TAL23 Other** VPH: Full or Ranges only TEPH EPH: Full or Ranges only ME4215 TPH: 8015 (Diesel Range) ME4217 TPH: 8015 (Gas Range) ME4217 PCB: 8082 608 Soxhlet Y or N Pesticides: 8081 608 SVOC: 8260 524.2 624 VOC: 8260 524.2 624 Field Filtered Y or N		Matrix Key: C = Concrete WP = Wipe WW = Wastewater SW = Surface Water E = Extract		No. of Containers No. of pH checked Analytics Sample #	
Company: JOHN CRESSY		Address: 940 MAIN ST. LEWISTON ME 04240 Phone: (207) 765-5500 Quote #: PO# (if required): 12-32591		Sample Date Time		CB: 1050 N CB: 1102 N CB: 1106 N CB: 1115 N CB: 1121 N		CB: 1 CB: 1 CB: 1 CB: 1 CB: 1	
Comments, Additional Analyses, or Special Instructions:									
Email Results to: JOHN CRESSY									
Turnaround Time (TAT)		1 Day* 3 Days* 5 Days		2 Days* 4 Days* Standard (8-10 business days)		Level II* Level III* Level IV*		State: NH MA ME CT RI Other:	
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.									
Relinquished By Sampler: John Cressy		Date: 12-20 Date:		Time: 1/13/1 Time:		Received By: V Stromus Received By:			
Relinquished By: John Cressy		Date:		Time:		Received By:			



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 74875
CLIENT: Summit
PROJECT: Maine Energy

COOLER NUMBER: 74,223
NUMBER OF COOLERS: 2

A: PRELIMINARY EXAMINATION:

1. Cooler received by(initials): KB

DATE COOLER RECEIVED/OPENED: 02/20/13

2. Circle one:

Hand delivered
(If no, skip 3)

Shipped

3. Did cooler come with a shipping slip?

Y N

3a. Enter carrier name and airbill number here:

4. Were custody seals on the outside of cooler?

How many & where: _____ Seal Date: _____

Y N Seal Name: _____

5. Did the custody seals arrive unbroken and intact upon arrival?

Y N

6. COC#:

7. Were Custody papers filled out properly (ink,signed, legible, project information etc)?

Y N

8. Were custody papers sealed in a plastic bag?

Y N

9. Did you sign the COC in the appropriate place?

Y N

10. Was enough ice used to chill the cooler?

Y N Temp. of cooler:

0.3°C, 0.7°C

B. Log-In: Date samples were logged in:

By: KB 02/20/13

11. Were all bottles sealed in separate plastic bags?

Y N

12. Did all bottles arrive unbroken and were labels in good condition?

Y N

13. Were all bottle labels complete(ID,Date,time,etc.)

Y N

14. Did all bottle labels agree with custody papers?

Y N

15. Were the correct containers used for the tests indicated:

Y N

16. Were samples received at the correct pH?

Y N

17. Was sufficient amount of sample sent for the tests indicated?

Y N

18. Were all samples submitted within holding time?

Y N

19. Were all containers used within AEL's expiration date?**

Y N

20. Were VOA samples absent of greater than pea-sized bubbles?

(Note:Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

Y N*

*If NO, List Sample ID's, Lab #s: _____

When bubbles are present in VOA samples they are labelled from smallest (no bubbles) to largest. Lab to analyze VOA samples with no bubbles or
smallest bubbles first

20. Laboratory labeling verified by (initials): JB

Date: 2/20/13

**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.

March 1, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74886
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

- Case Narrative/Non-Conformance Summary
- Sample Log Sheet - Cover Page
- PCB Form 1 Data Sheet for Samples
- PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
 - Chromatograms
- Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No exceptions.

PCBs by EPA Method 8082:
No results were reported below the quantitation limit.

All samples except 74886-13 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument or matrix affect.

The MS/MSD analyzed on sample 74886-1 did not meet acceptance criteria for PCB 1260 recoveries due to the parent sample having concentrations of PCB 1260 that exceeded the calibration range of the instrument. The laboratory control samples (L022513PSOX/LD022513PSOX) were in control for all recoveries and RPDs. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC



Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
Summit Environmental Consultants Inc.
640 Main Street
Lewiston ME 04240

Report Number: 74886

Revision: Rev. 0

Re: MAINE ENERGY (Project No: 12-3259.1)

Enclosed are the results of the analyses on your sample(s). Samples were received on 22 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Sample Analysis: The attached pages detail the Client Sample IDs, Lab Sample IDs, and Analyses requested

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

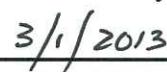
If you have any questions on these results, please do not hesitate to contact us.

Authorized signature



Stephen L. Knollmeyer Lab. Director

Date



**This report shall not be reproduced, except in full, without the written
consent of Analytics Environmental Laboratory, LLC.**

CLIENT: Summit Environmental
 Consultants Inc.

REPORT NUMBER: 74886

REV: Rev. 0

PROJECT: MAINE ENERGY (Project No: 12-3259.1)

<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74886-1	02/20/13	SE-SB-234 (2-4')	EPA 8082 (PCBs only)	
74886-2	02/20/13	SE-SB-234 (4-6')	EPA 8082 (PCBs only)	
74886-3	02/20/13	SE-SB-234 (6-8')	EPA 8082 (PCBs only)	
74886-4	02/20/13	SE-SB-234 (8-10')	EPA 8082 (PCBs only)	
74886-5	02/20/13	SE-SB-230 (0-2')	EPA 8082 (PCBs only)	
74886-6	02/20/13	SE-SB-230 (2-4')	EPA 8082 (PCBs only)	
74886-7	02/20/13	SE-SB-231 (0-2')	EPA 8082 (PCBs only)	
74886-8	02/20/13	SE-SB-231 (2-4')	EPA 8082 (PCBs only)	
74886-9	02/20/13	SE-SB-231 (4-6')	EPA 8082 (PCBs only)	
74886-10	02/20/13	SE-SB-231 (6-8')	EPA 8082 (PCBs only)	
74886-11	02/20/13	SE-SB-235 (0-2')	EPA 8082 (PCBs only)	
74886-12	02/20/13	SE-SB-235 (2-4')	EPA 8082 (PCBs only)	
74886-13	02/20/13	SE-SB-235 (4-6')	EPA 8082 (PCBs only)	
74886-14	02/20/13	SE-SB-235 (6-8')	EPA 8082 (PCBs only)	
74886-15	02/20/13	SE-SB-235 (8-10')	EPA 8082 (PCBs only)	
74886-16	02/20/13	SE-SB-229 (0-2')	EPA 8082 (PCBs only)	
74886-17	02/20/13	SE-SB-229 (2-4')	EPA 8082 (PCBs only)	
74886-18	02/20/13	SE-SB-229 (4-6')	EPA 8082 (PCBs only)	
74886-19	02/20/13	SE-SB-229 (6-8')	EPA 8082 (PCBs only)	
74886-20	02/20/13	SE-SB-229 (8-10')	EPA 8082 (PCBs only)	

Surrogate Compound Limits

	Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water				
1,4-Difluorobenzene		70-130		EPA 524.2
Bromofluorobenzene		70-130		
1,2-Dichlorobenzene-d4		70-130		
Volatile Organic Compounds				
1,2-Dichloroethane-d4		70-120	70-120	EPA 624/8260B
Toluene-d8		85-120	85-120	
Bromofluorobenzene		75-120	75-120	
Semi-Volatile Organic Compounds				
2-Fluorophenol		20-110	35-105	EPA 625/8270C
d5-Phenol		15-110	40-100	
d5-nitrobenzene		40-110	35-100	
2-Fluorobiphenyl		50-110	45-105	
2,4,6-Tribromophenol		40-110	40-125	
d14-p-terphenyl		50-130	30-125	
PAH's by SIM				
d5-nitrobenzene		21-110	35-110	EPA 8270C
2-Fluorobiphenyl		36-121	45-105	
d14-p-terphenyl		33-141	30-125	
Pesticides and PCBs				
2,4,5,6-Tetrachloro-m-xylene (TCX)		46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)		40-135	40-130	
Herbicides				
Dichloroacetic acid (DCAA)		30-150	30-150	
Gasoline Range Organics/TPH Gasoline				
Trifluorotoluene TFT (FID)		60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)		60-140	60-140	
Trifluorotoluene TFT (PID)		60-140	60-140	
Bromofluorobenzene (BFB) (PID)		60-140	60-140	
Diesel Range Organics/TPH Diesel				
m-terphenyl		60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons				
2,5-Dibromotoluene (PID)		70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)		70-130	70-130	
Extricable Petroleum Hydrocarbons				
1-chloro-octadecane (aliphatic)		40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)		40-140	40-140	
2-Fluorobiphenyl (Fractionation)		40-140	40-140	
2-Bromonaphthalene (fractionation)		40-140	40-140	

PCB
DATA SUMMARIES

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-234 (2-4')

Lab Sample ID: 74886-1
 Matrix: Solid
 Percent Solid: 89
 Dilution Factor: 11
 Collection Date: 02/20/13
 Lab Receipt Date: 02/22/13
 Extraction Date: 02/22/13
 Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	2350

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	85	%
Decachlorobiphenyl	82	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-1,1:10,A/C

Column ID: 0.25 mm

Data File: M67690.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2345	2354	0.4	

Column to be used to flag RPD values greater than QC limit of 40%

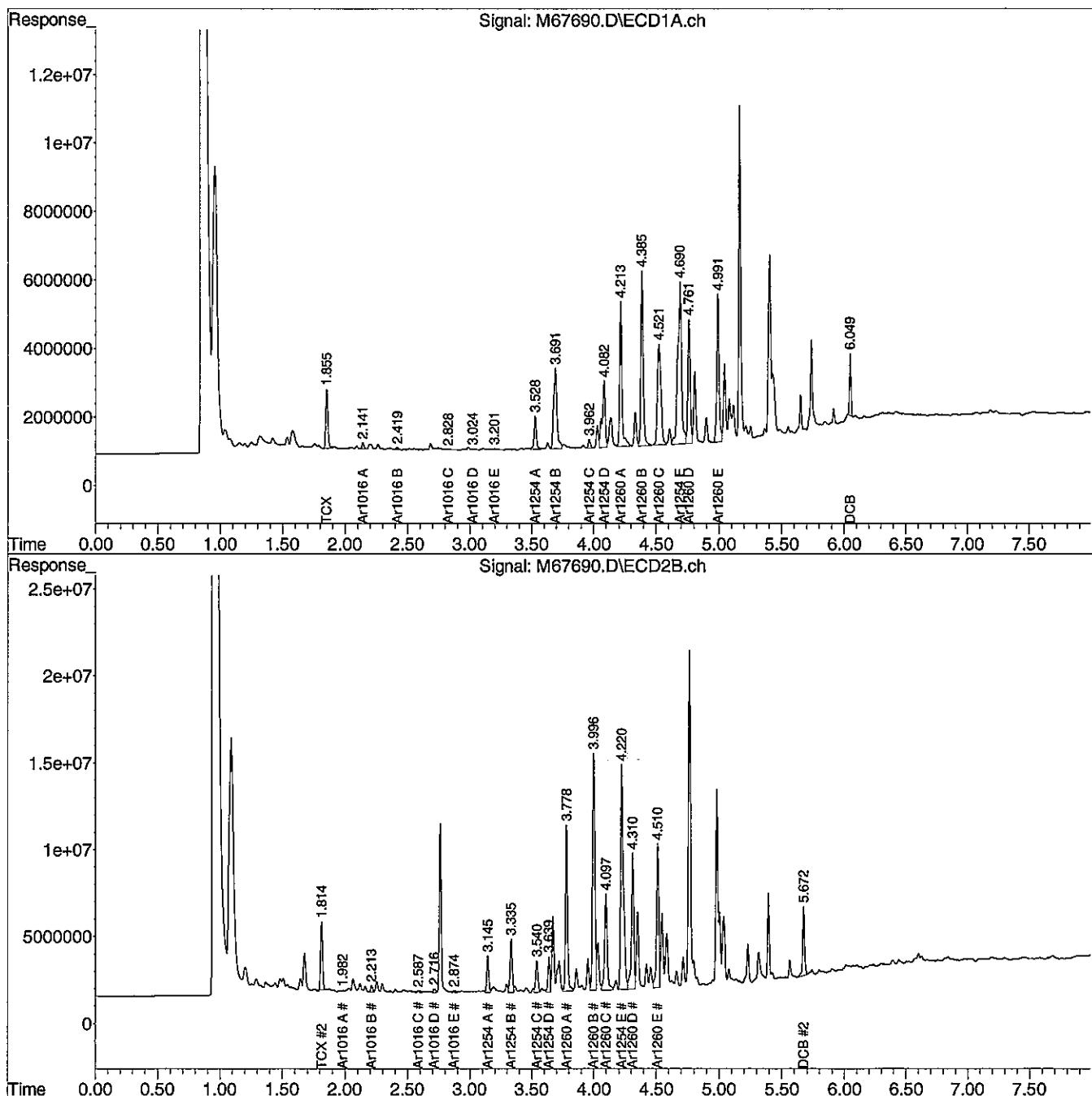
* Values outside QC limits

Comments: _____

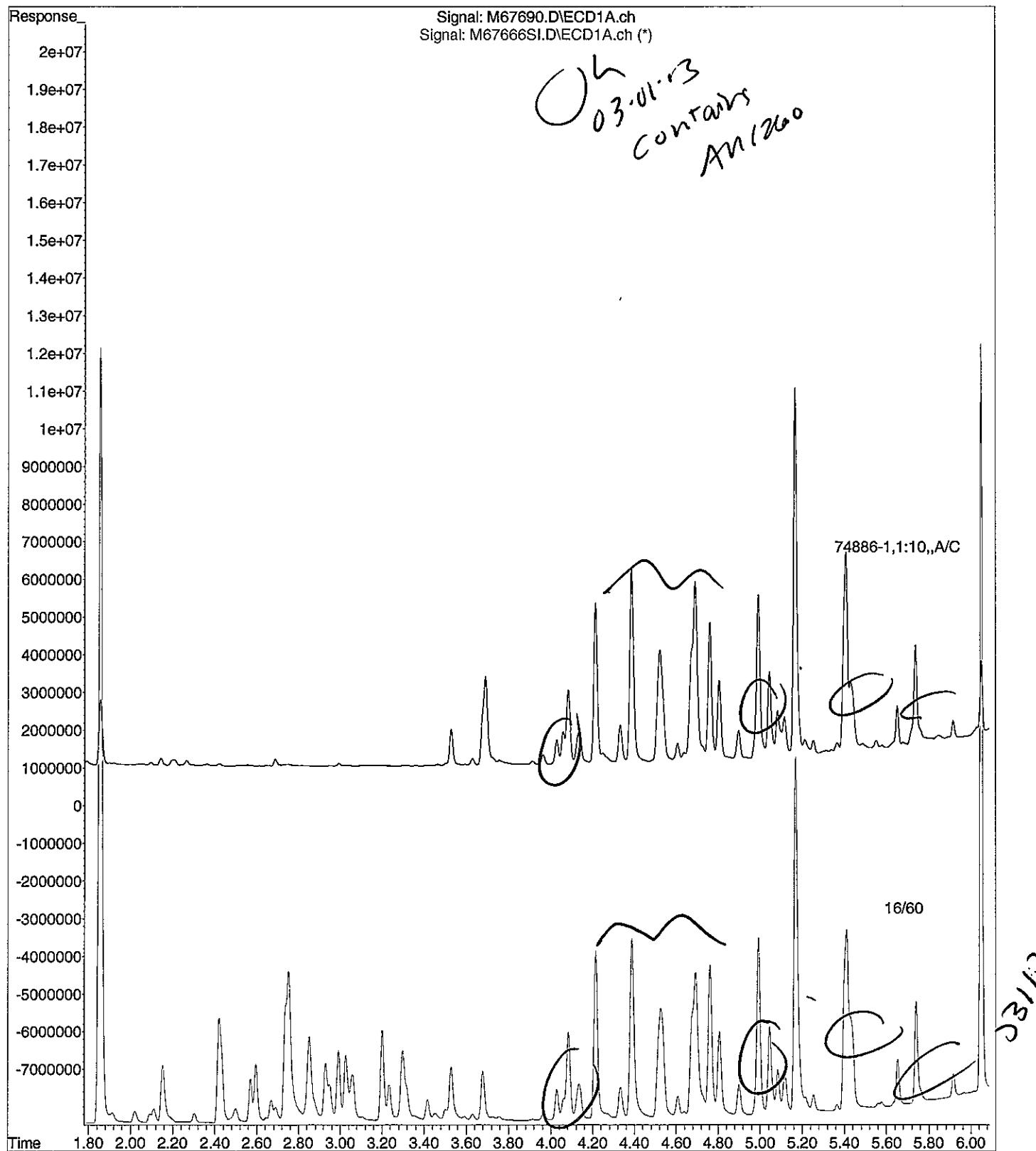
Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67690.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 10:40 pm
 Operator : JK
 Sample : 74886-1,1:10,,A/C
 Misc : SOIL
 ALS Vial : 34 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:19 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022813-M\M67690.D
Operator : JK
Acquired : 28 Feb 2013 10:40 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74886-1,1:10,,A/C
Misc Info : SOIL
Vial Number: 34



Mr. John Cressey
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 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-234 (4-6")

Lab Sample ID: 74886-2
Matrix: Solid
Percent Solid: 89
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	1840

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-2,1:10,,A/C

Column ID: 0.25 mm

Data File: M67693.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1775	1840	3.6	

Column to be used to flag RPD values greater than QC limit of 40%

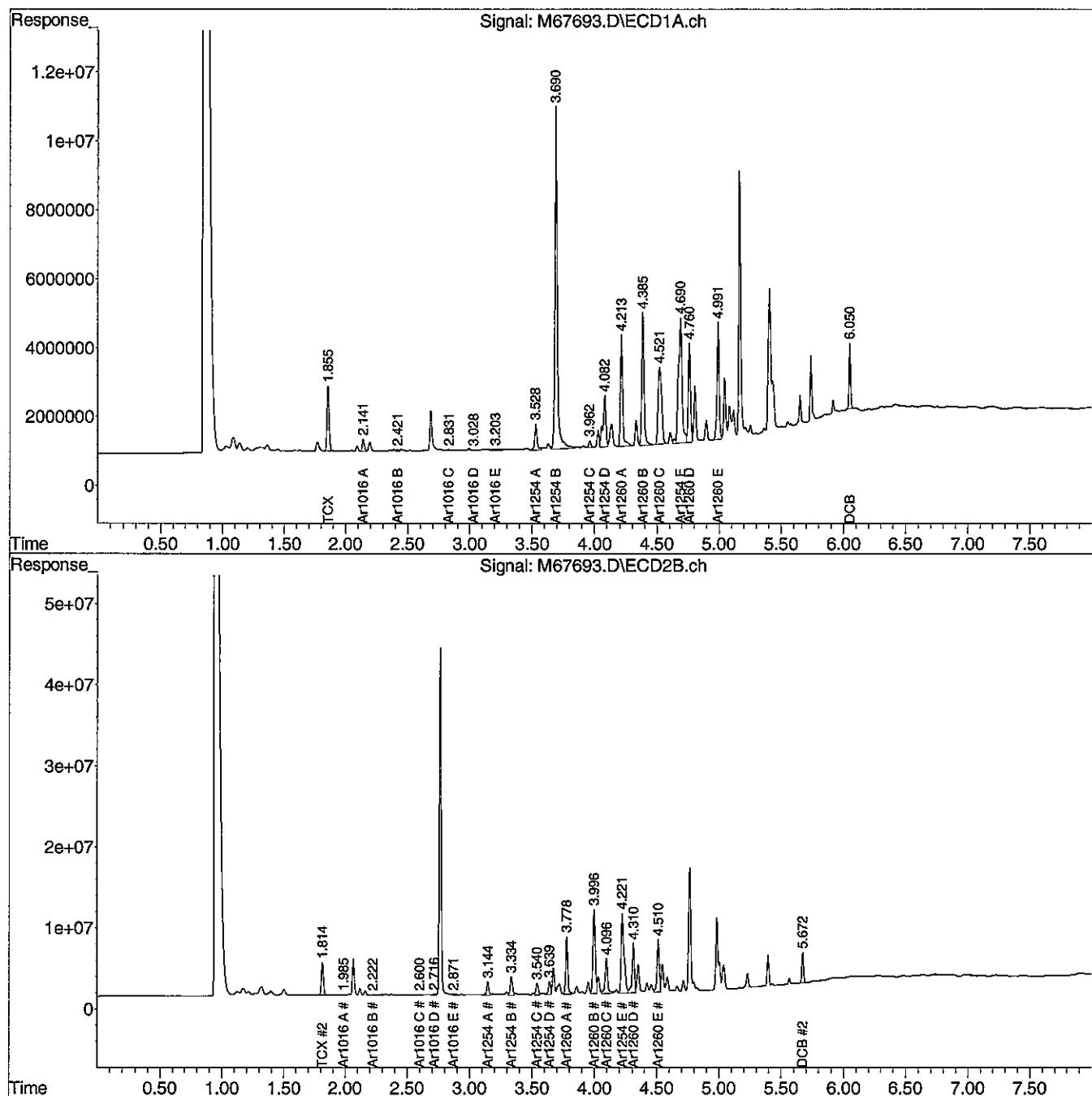
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67693.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:11 pm
 Operator : JK
 Sample : 74886-2,1:10,,A/C
 Misc : SOIL
 ALS Vial : 37 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:25 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 640 Main Street
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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-234 (6-8")

Lab Sample ID: 74886-3
Matrix: Solid
Percent Solid: 86
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	2260

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	94	%
Decachlorobiphenyl	94	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-3,1:10,,A/C

Column ID: 0.25 mm

Data File: M67694.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 11.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2250	2255	0.2	

Column to be used to flag RPD values greater than QC limit of 40%

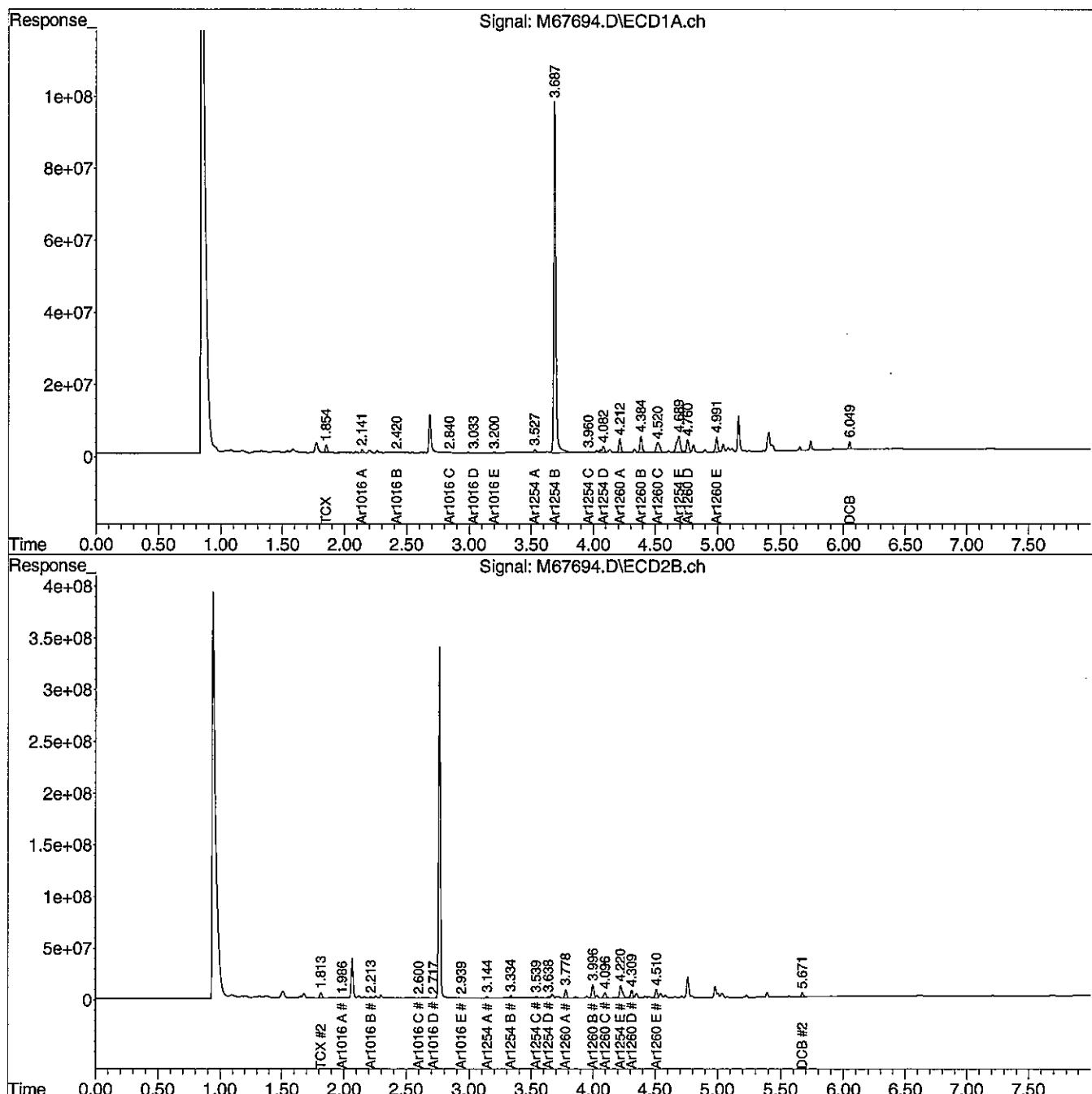
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67694.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:21 pm
 Operator : JK
 Sample : 74886-3,1:10,,A/C
 Misc : SOIL
 ALS Vial : 38 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:27 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-234 (8-10')

Lab Sample ID: 74886-4
Matrix: Solid
Percent Solid: 90
Dilution Factor: 10
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	330	U
PCB-1221	330	U
PCB-1232	330	U
PCB-1242	330	U
PCB-1248	330	U
PCB-1254	330	U
PCB-1260	330	1540

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	88	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-4,1:10,,A/C

Column ID: 0.25 mm

Data File: M67695.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1535	1533	0.1	

Column to be used to flag RPD values greater than QC limit of 40%

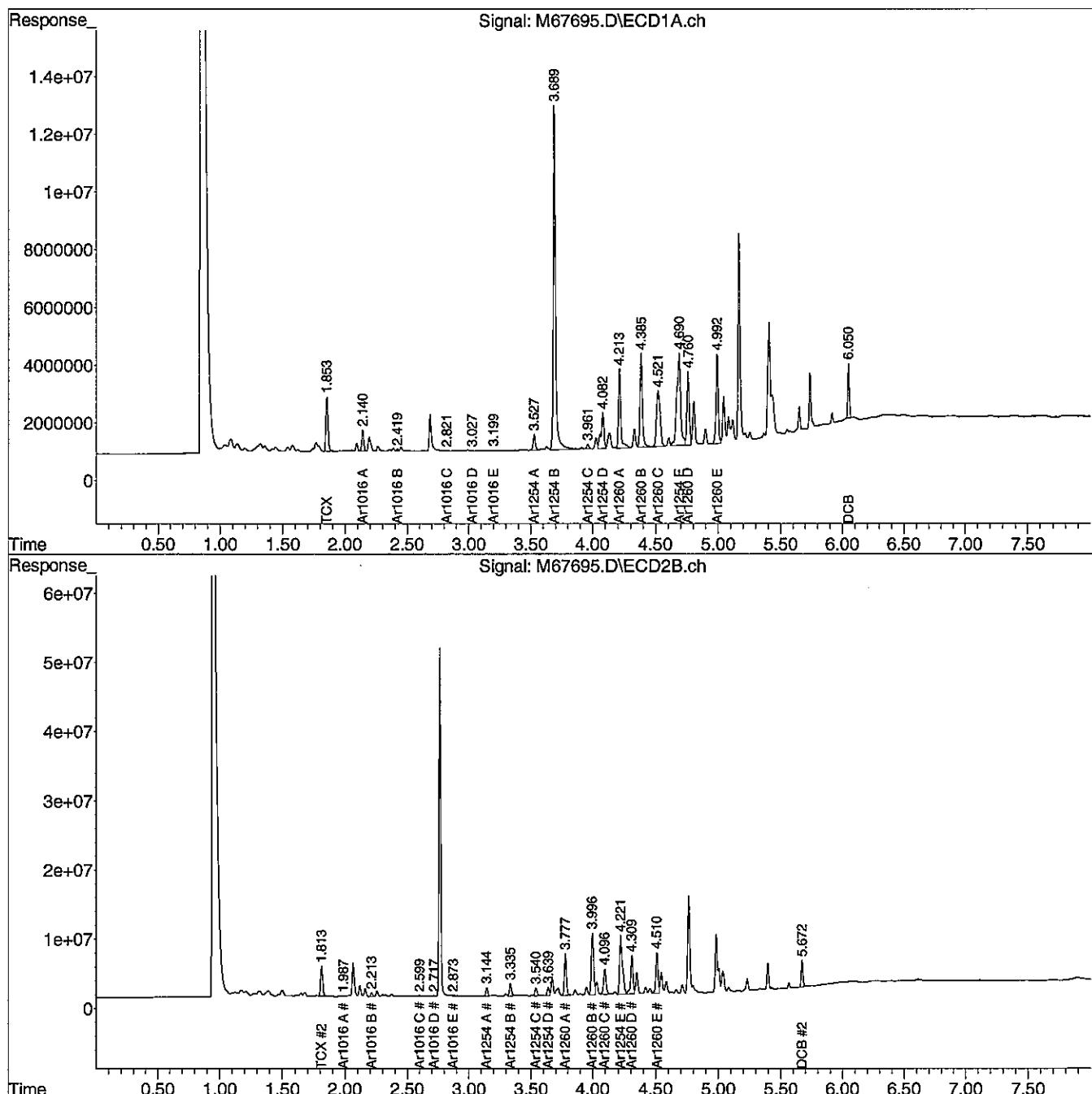
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67695.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:31 pm
 Operator : JK
 Sample : 74886-4,1:10,,A/C
 Misc : SOIL
 ALS Vial : 39 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:29 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 640 Main Street
 Lewiston ME 04240

March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-230 (0-2")

Lab Sample ID: 74886-5
 Matrix: Solid
 Percent Solid: 80
 Dilution Factor: 115
 Collection Date: 02/20/13
 Lab Receipt Date: 02/22/13
 Extraction Date: 02/22/13
 Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	3800	U
PCB-1221	3800	U
PCB-1232	3800	U
PCB-1242	3800	U
PCB-1248	3800	U
PCB-1254	3800	U
PCB-1260	3800	27700

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-5,1:100,,A/C

Column ID: 0.25 mm

Data File: M67696.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 114.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	27689	24630	11.7	

Column to be used to flag RPD values greater than QC limit of 40%

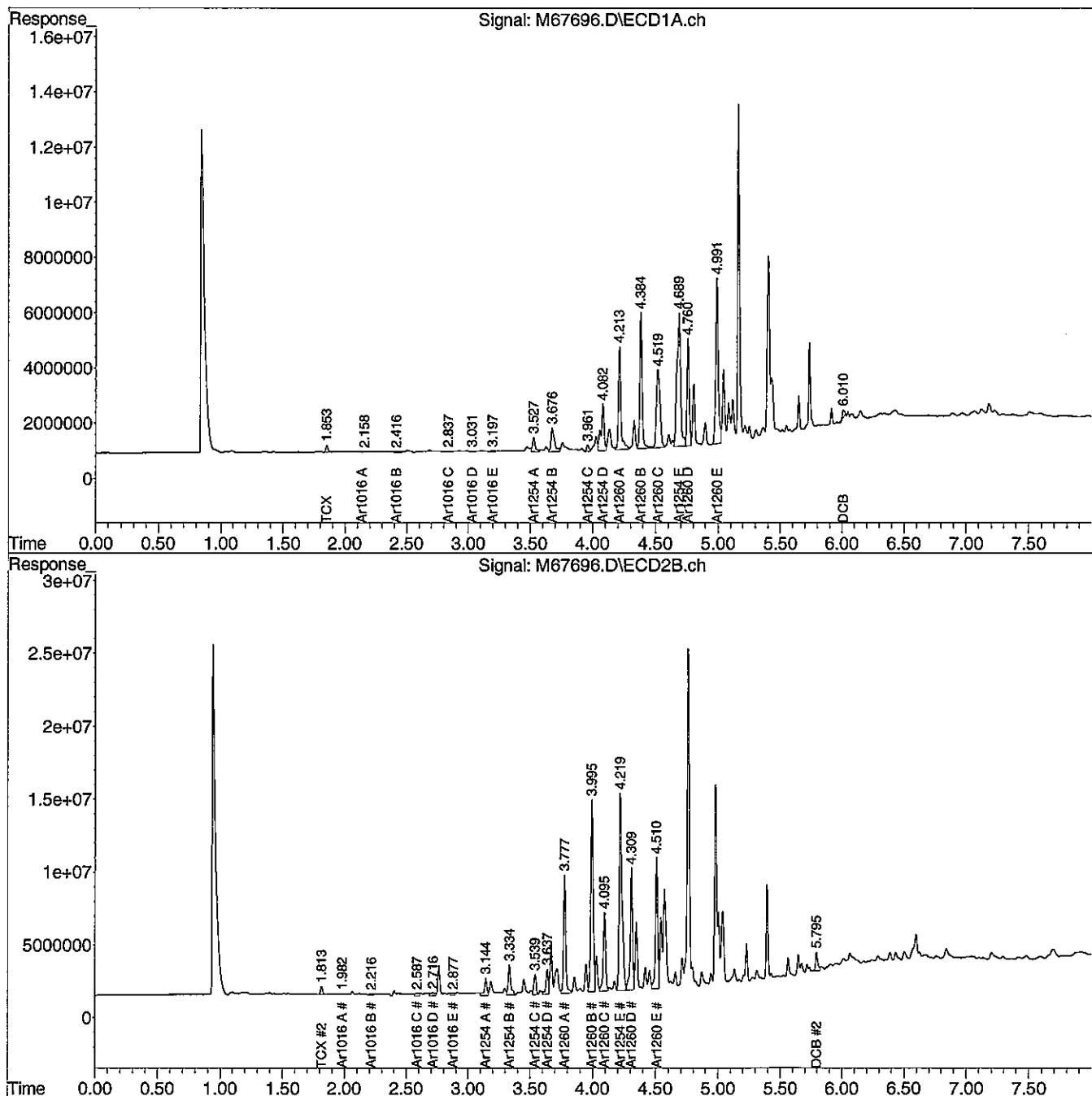
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67696.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:41 pm
 Operator : JK
 Sample : 74886-5, 1:100, , A/C
 Misc : SOIL
 ALS Vial : 40 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:31 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-230 (2-4')

Lab Sample ID: 74886-6
Matrix: Solid
Percent Solid: 92
Dilution Factor: 10
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	330	U
PCB-1221	330	U
PCB-1232	330	U
PCB-1242	330	U
PCB-1248	330	U
PCB-1254	330	U
PCB-1260	330	3080

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	103	%	
Decachlorobiphenyl	95	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-6,1:10,,A/C

Column ID: 0.25 mm

Data File: M67697.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3081	3015	2.1	

Column to be used to flag RPD values greater than QC limit of 40%

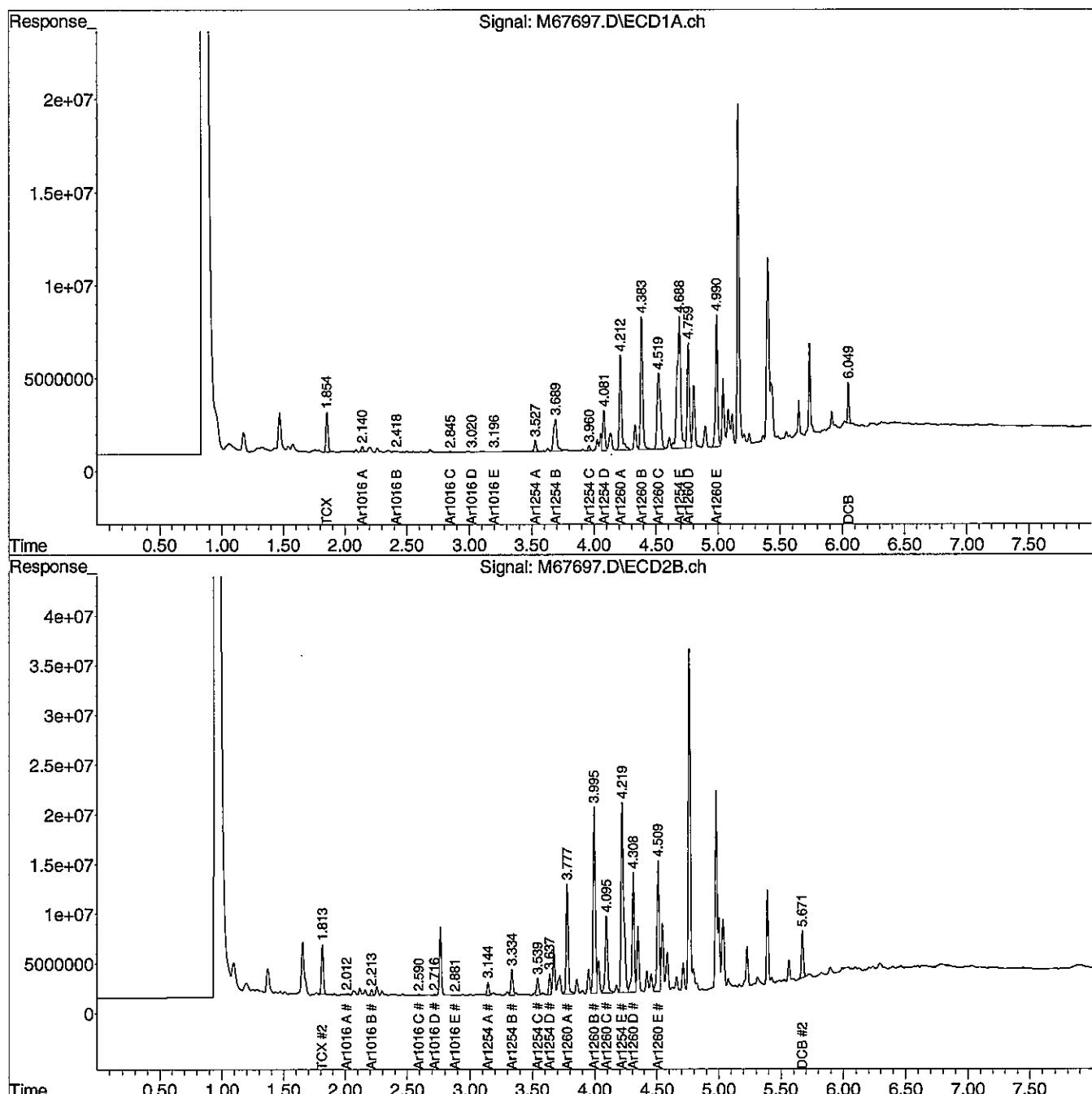
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67697.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:51 pm
 Operator : JK
 Sample : 74886-6,1:10,,A/C
 Misc : SOIL
 ALS Vial : 41 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:33 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



Mr. John Cressey
 Summit Environmental Consultants Inc.
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 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-231 (0-2')

Lab Sample ID: 74886-7
Matrix: Solid
Percent Solid: 86
Dilution Factor: 56
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	1850	U
PCB-1221	1850	U
PCB-1232	1850	U
PCB-1242	1850	U
PCB-1248	1850	U
PCB-1254	1850	U
PCB-1260	1850	17800

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-7,1:50,,A/C

Column ID: 0.25 mm

Data File: M67698.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 56.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	17767	17761	0.0	

Column to be used to flag RPD values greater than QC limit of 40%

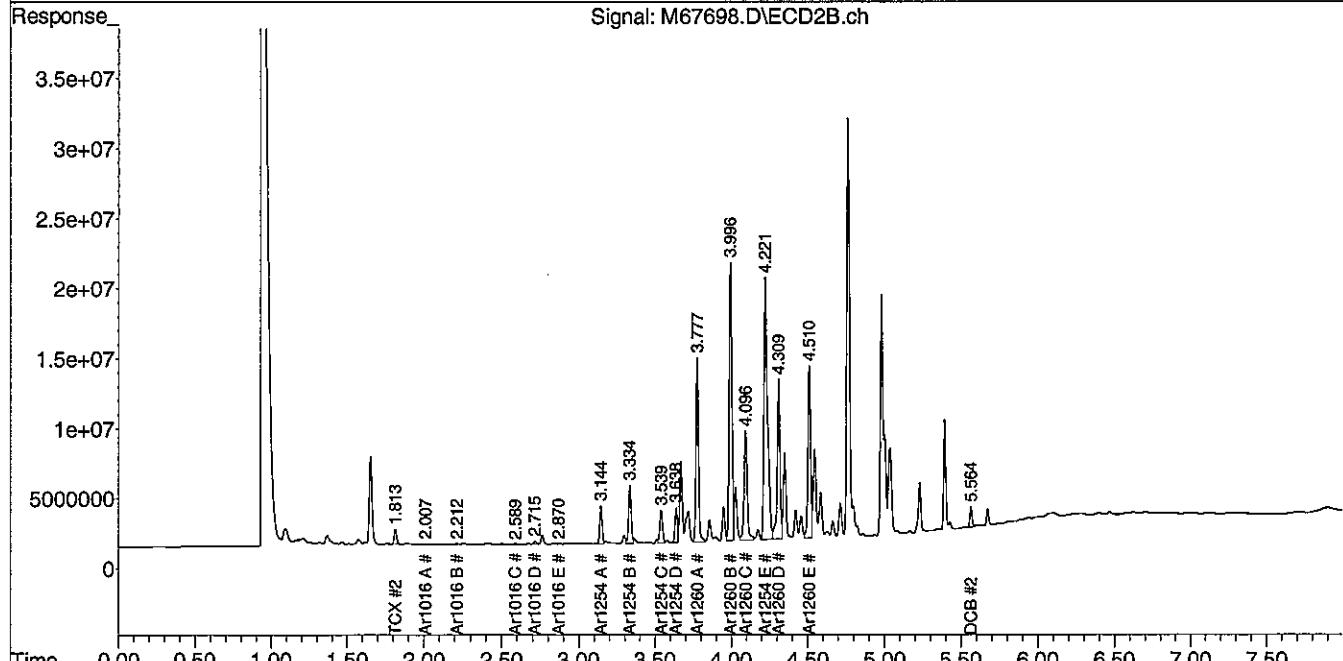
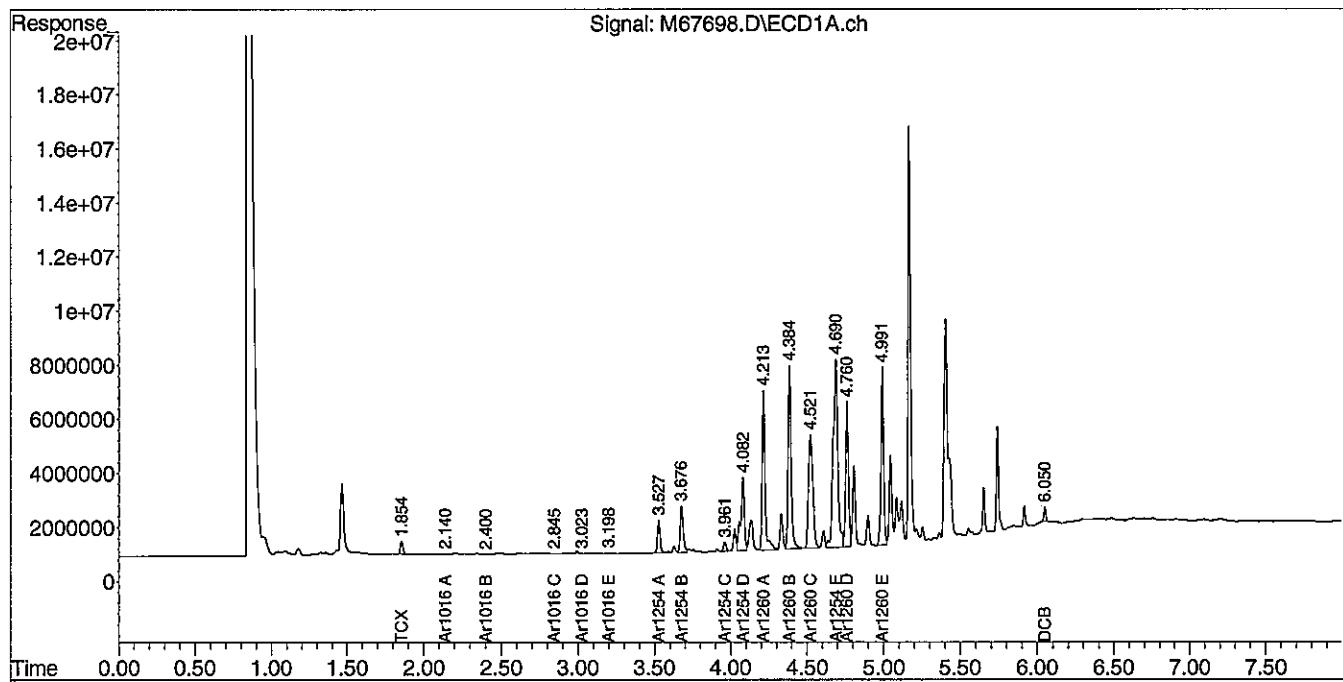
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67698.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:01 am
 Operator : JK
 Sample : 74886-7,1:50,,A/C
 Misc : SOIL
 ALS Vial : 42 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:35 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-231 (2-4")

Lab Sample ID: 74886-8
Matrix: Solid
Percent Solid: 89
Dilution Factor: 1090
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	36000	U
PCB-1221	36000	U
PCB-1232	36000	U
PCB-1242	36000	U
PCB-1248	36000	U
PCB-1254	36000	U
PCB-1260	36000	324000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-8,1:1000,,A/C

Column ID: 0.25 mm

Data File: M67699.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1090.3

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	321505	324249	0.8	

Column to be used to flag RPD values greater than QC limit of 40%

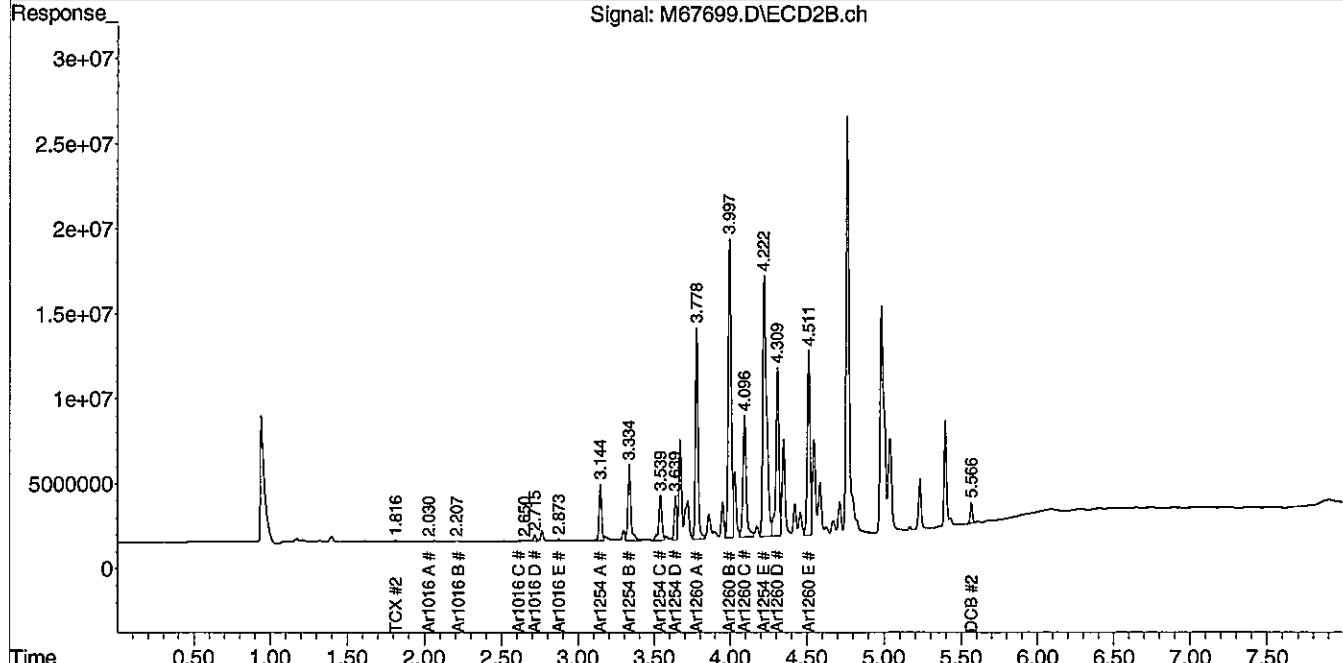
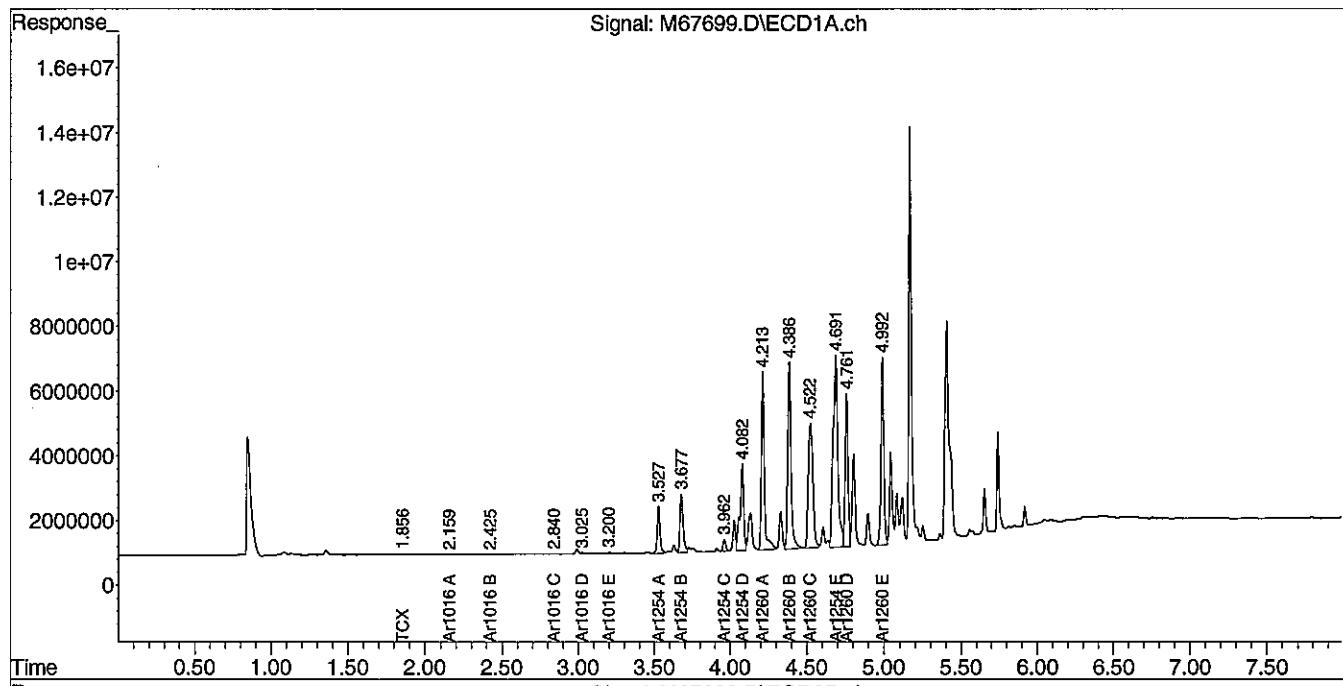
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67699.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:11 am
 Operator : JK
 Sample : 74886-8,1:1000,,A/C
 Misc : SOIL
 ALS Vial : 43 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:37 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-231 (4-6')

Lab Sample ID:	74886-9
Matrix:	Solid
Percent Solid:	89
Dilution Factor:	217
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/22/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	7160	U
PCB-1221	7160	U
PCB-1232	7160	U
PCB-1242	7160	U
PCB-1248	7160	U
PCB-1254	7160	U
PCB-1260	7160	164000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-9,1:200,,A/C

Column ID: 0.25 mm

Data File: M67700.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 217.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	163775	163769	0.0	

Column to be used to flag RPD values greater than QC limit of 40%

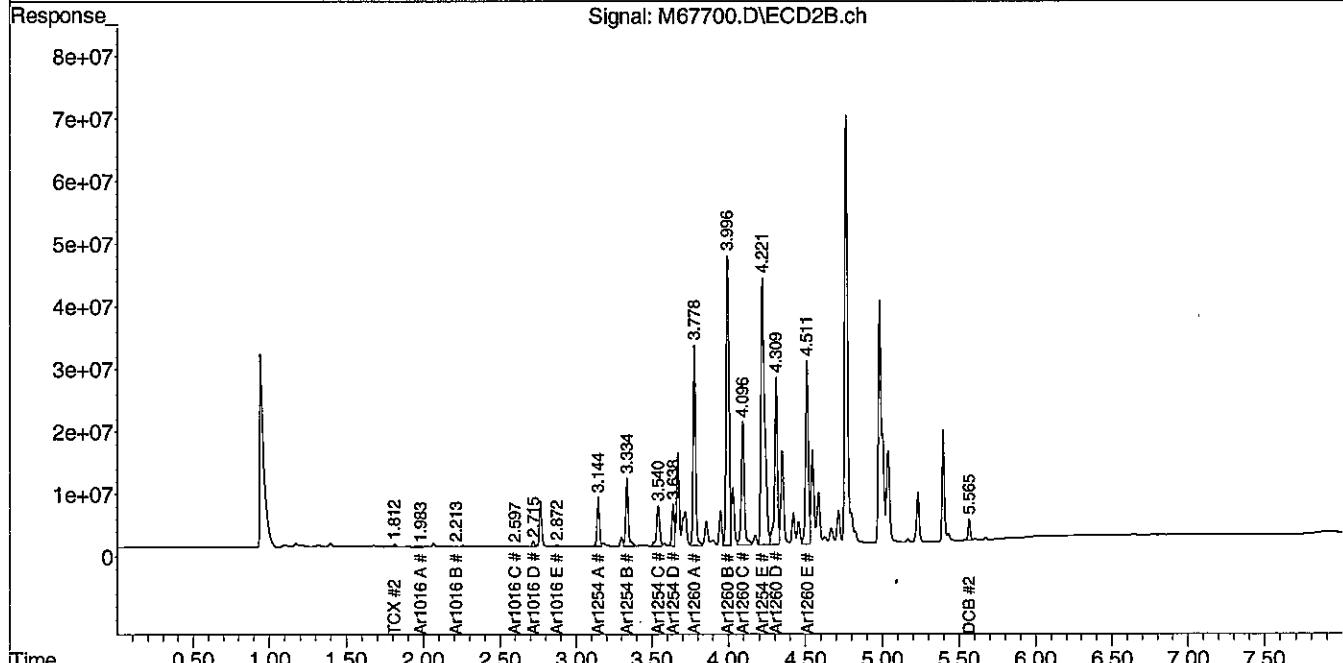
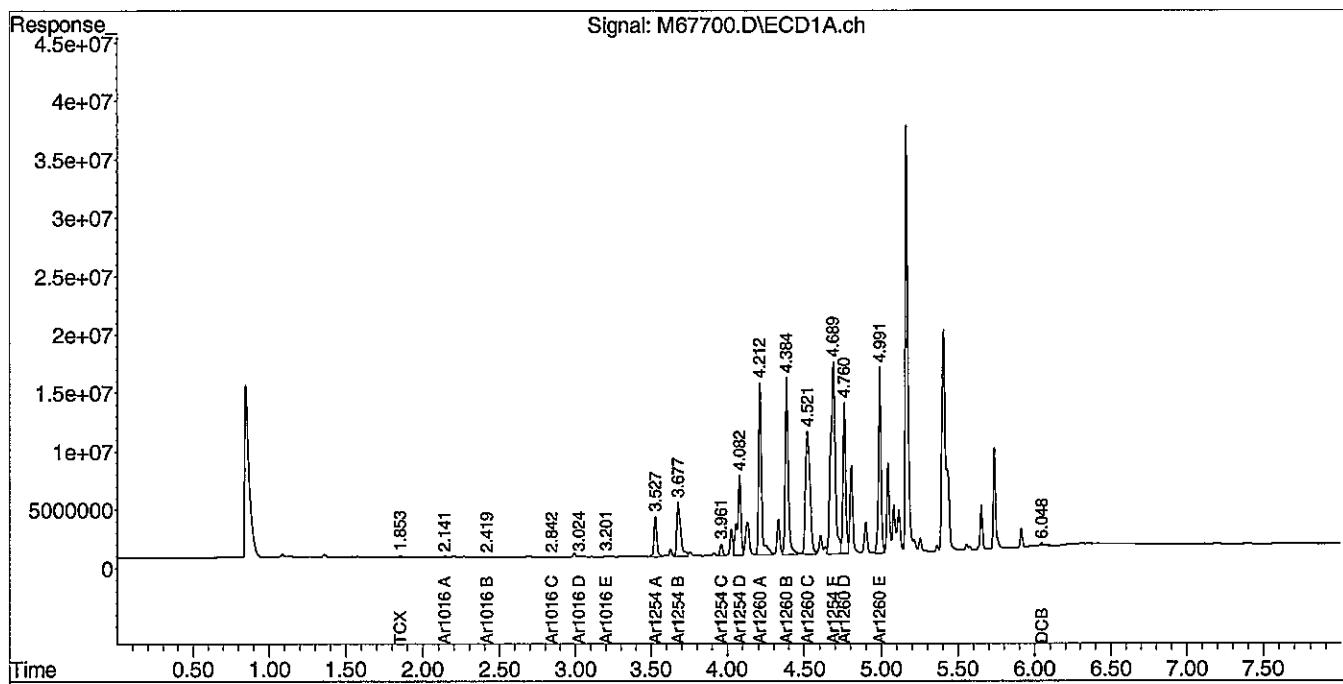
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67700.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:21 am
 Operator : JK
 Sample : 74886-9,1:200,,A/C
 Misc : SOIL
 ALS Vial : 44 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:39 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-231 (6-8")

Lab Sample ID: 74886-10
Matrix: Solid
Percent Solid: 89
Dilution Factor: 5320
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/22/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	176000	U
PCB-1221	176000	U
PCB-1232	176000	U
PCB-1242	176000	U
PCB-1248	176000	U
PCB-1254	176000	U
PCB-1260	176000	3150000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-10,1:5000,,A/C

Column ID: 0.25 mm

Data File: M67767.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5319.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3153212	2907949	8.1	

Column to be used to flag RPD values greater than QC limit of 40%

* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\023013-M\
 Data File : M67767.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:38 pm
 Operator : JK
 Sample : 74886-10,1:5000,,A/C
 Misc : SOIL
 ALS Vial : 10 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e

Quant Time: Mar 01 15:39:40 2013

Quant Method : C:\msdchem\1\METHODS\PCB020513.M

Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254

QLast Update : Thu Feb 28 11:22:13 2013

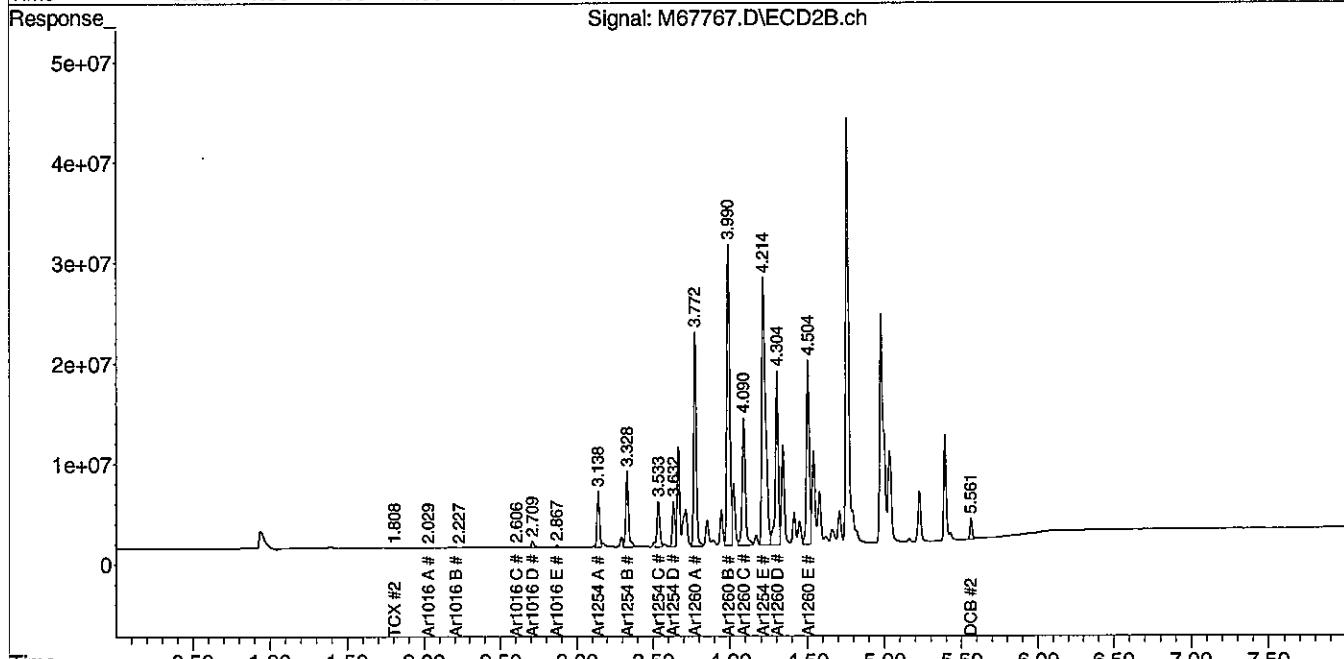
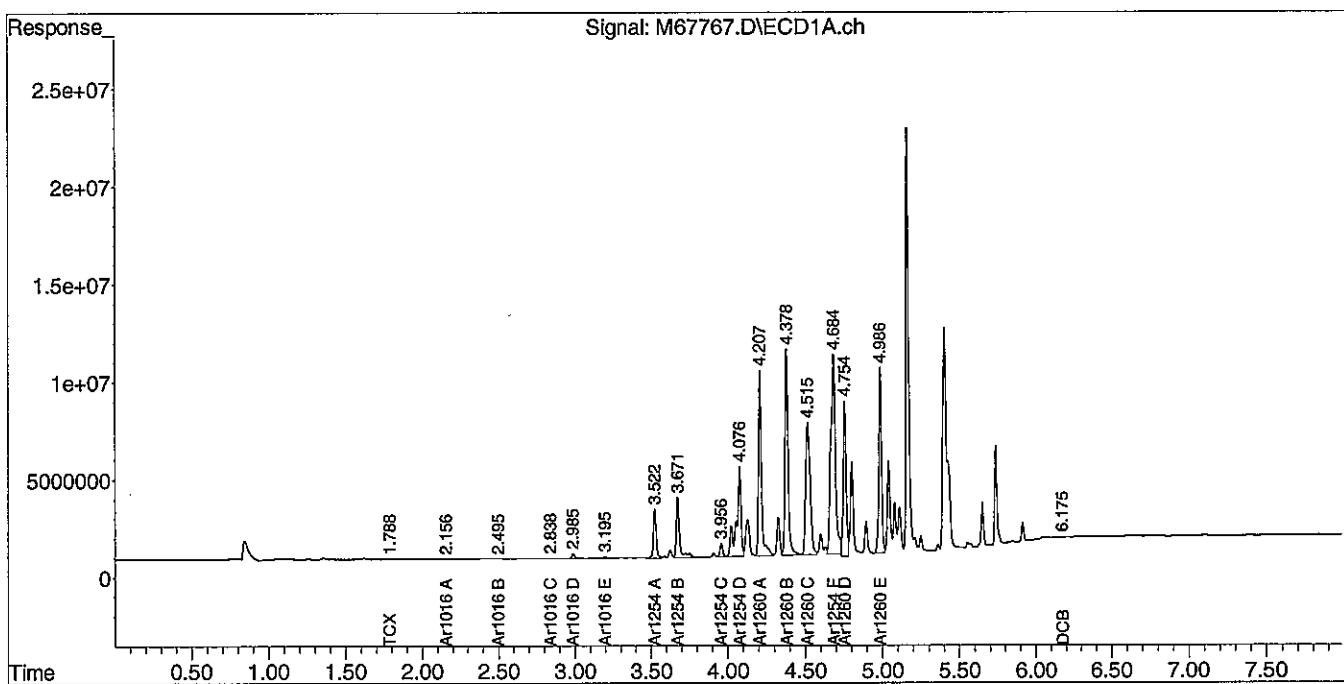
Response via : Initial Calibration

Integrator: ChemStation

Volume Inj. : 2 uL

Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides

Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-235 (0-2')

Lab Sample ID: 74886-11
Matrix: Solid
Percent Solid: 89
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	2240

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	73	%
Decachlorobiphenyl	67	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-11,1:5

Column ID: 0.25 mm

Data File: M67702.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.5

Column ID: 0.25 mm

COMPOUND	Column #1		Column #2	
	SAMPLE RESULT (ug/kg)		SAMPLE RESULT (ug/kg)	RPD
PCB 1260	2240		2188	2.3

Column to be used to flag RPD values greater than QC limit of 40%

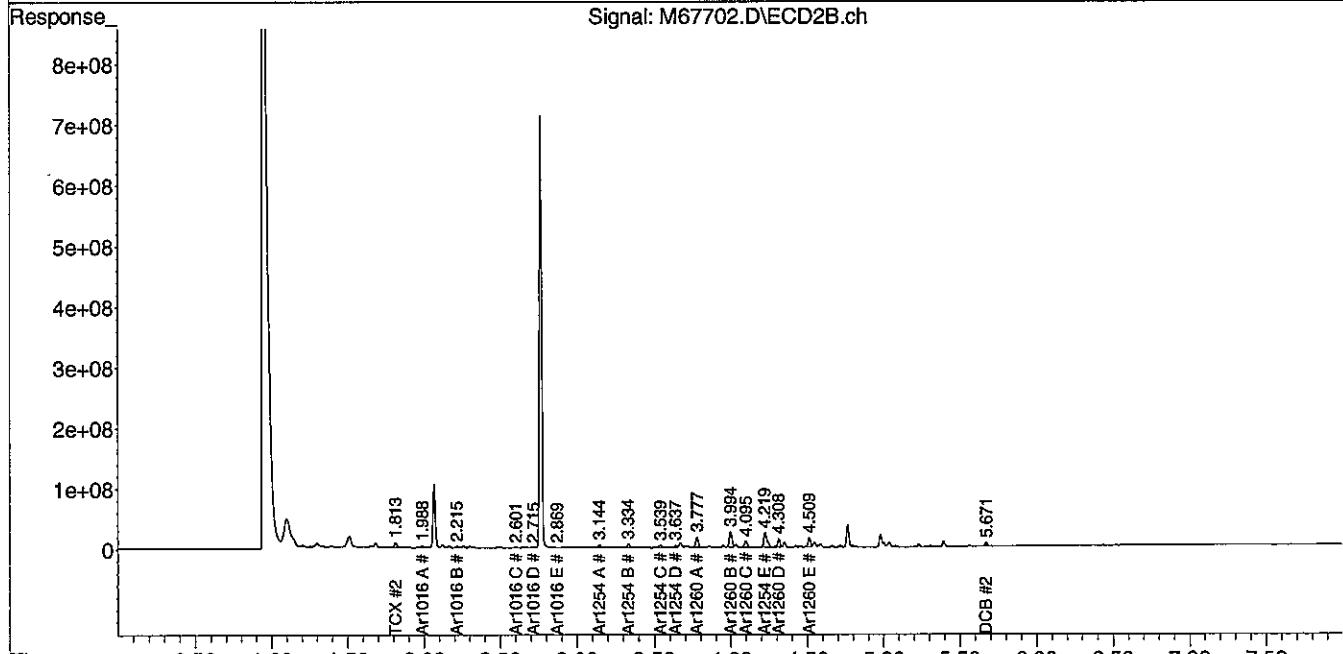
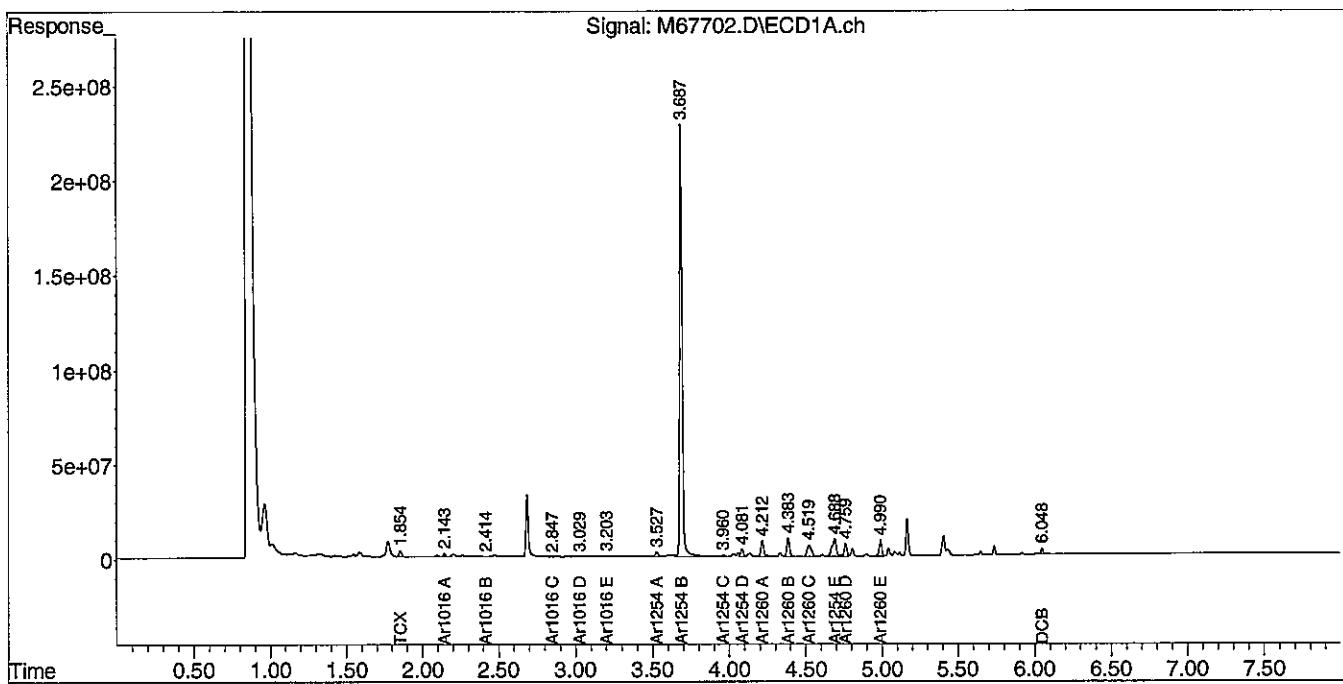
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67702.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:42 am
 Operator : JK
 Sample : 74886-11,1:5
 Misc : SOIL
 ALS Vial : 46 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:43 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-235 (2-4")

Lab Sample ID:	74886-12
Matrix:	Solid
Percent Solid:	87
Dilution Factor:	2.2
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	73	U
PCB-1221	73	U
PCB-1232	73	U
PCB-1242	73	U
PCB-1248	73	U
PCB-1254	73	U
PCB-1260	73	488

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	72	%
Decachlorobiphenyl	77	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-12.1:2

Column ID: 0.25 mm

Data File: M67703.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	488	440	10.4	

Column to be used to flag RPD values greater than QC limit of 40%

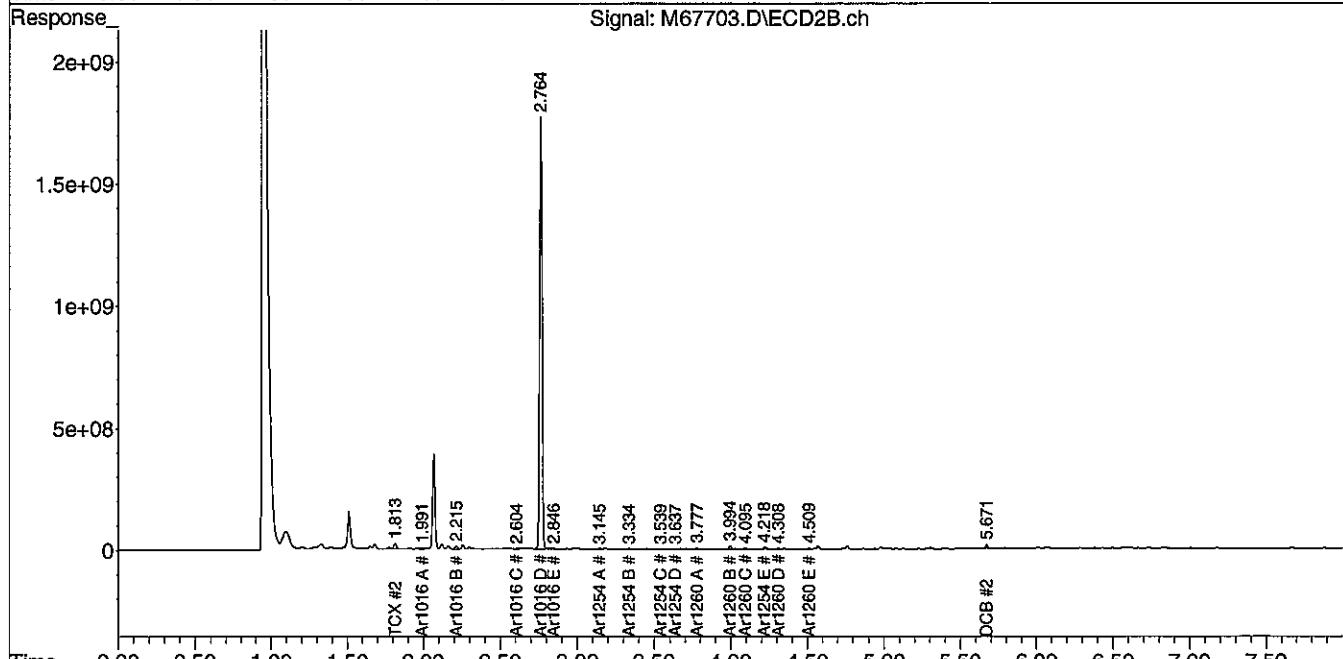
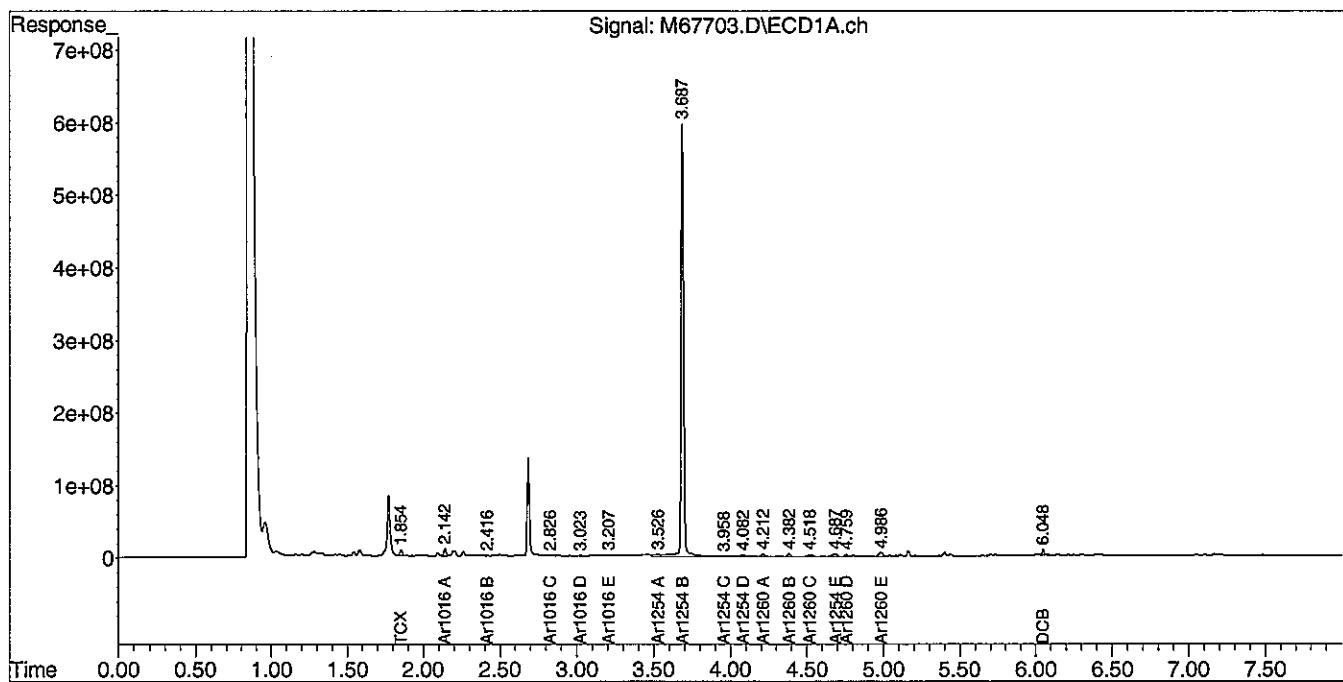
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67703.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 12:52 am
 Operator : JK
 Sample : 74886-12.1:2
 Misc : SOIL
 ALS Vial : 47 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:45 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-235 (4-6')

Lab Sample ID: 74886-13
Matrix: Solid
Percent Solid: 89
Dilution Factor: 1.1
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	36	U
PCB-1221	36	U
PCB-1232	36	U
PCB-1242	36	U
PCB-1248	36	U
PCB-1254	36	U
PCB-1260	36	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	64	%
Decachlorobiphenyl	76	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

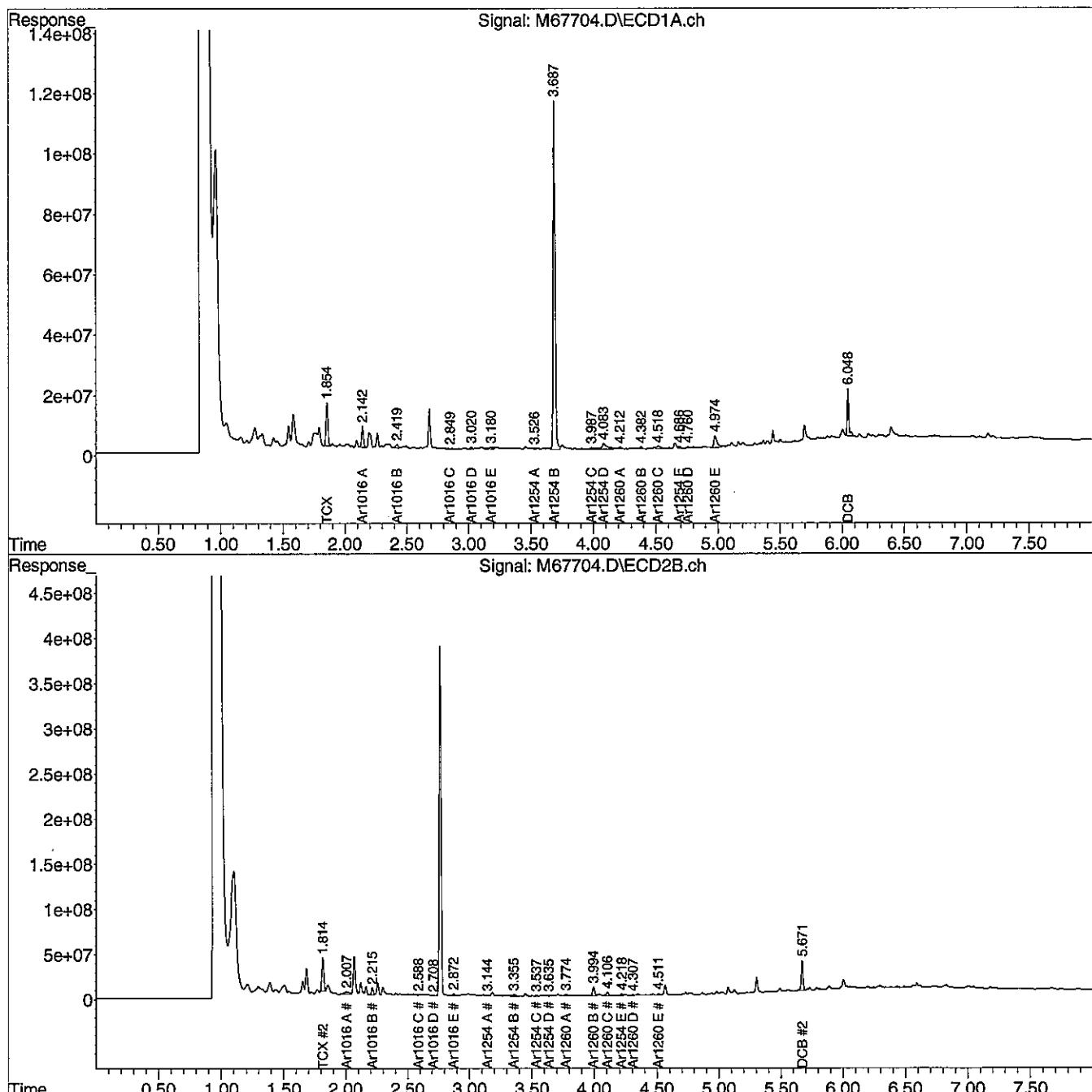
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67704.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 1:02 am
 Operator : JK
 Sample : 74886-13
 Misc : SOIL
 ALS Vial : 48 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:47 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-235 (6-8")

Lab Sample ID: 74886-14
Matrix: Solid
Percent Solid: 87
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	1380

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	80	%
Decachlorobiphenyl	72	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-14,1:5

Column ID: 0.25 mm

Data File: M67705.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.7

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1377	1353	1.7	

Column to be used to flag RPD values greater than QC limit of 40%

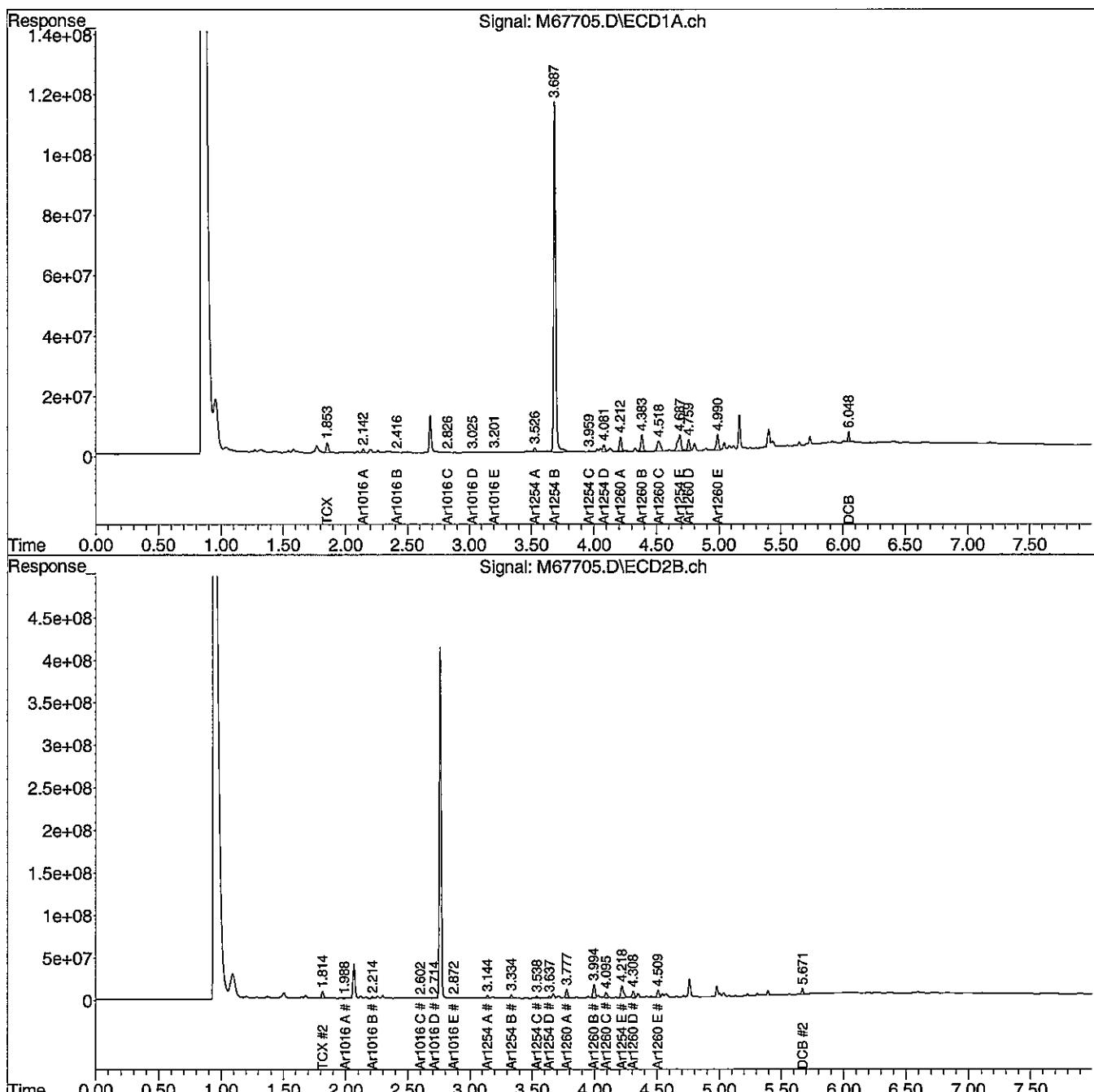
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67705.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 1:12 am
 Operator : JK
 Sample : 74886-14,1:5
 Misc : SOIL
 ALS Vial : 49 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:49 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-235 (8-10')

Lab Sample ID: 74886-15
Matrix: Solid
Percent Solid: 89
Dilution Factor: 2.2
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	73	U
PCB-1221	73	U
PCB-1232	73	U
PCB-1242	73	U
PCB-1248	73	U
PCB-1254	73	U
PCB-1260	73	399

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	68	%
Decachlorobiphenyl	65	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A.
 Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-15,1:2

Column ID: 0.25 mm

Data File: M67706.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	324	399	20.7	

Column to be used to flag RPD values greater than QC limit of 40%

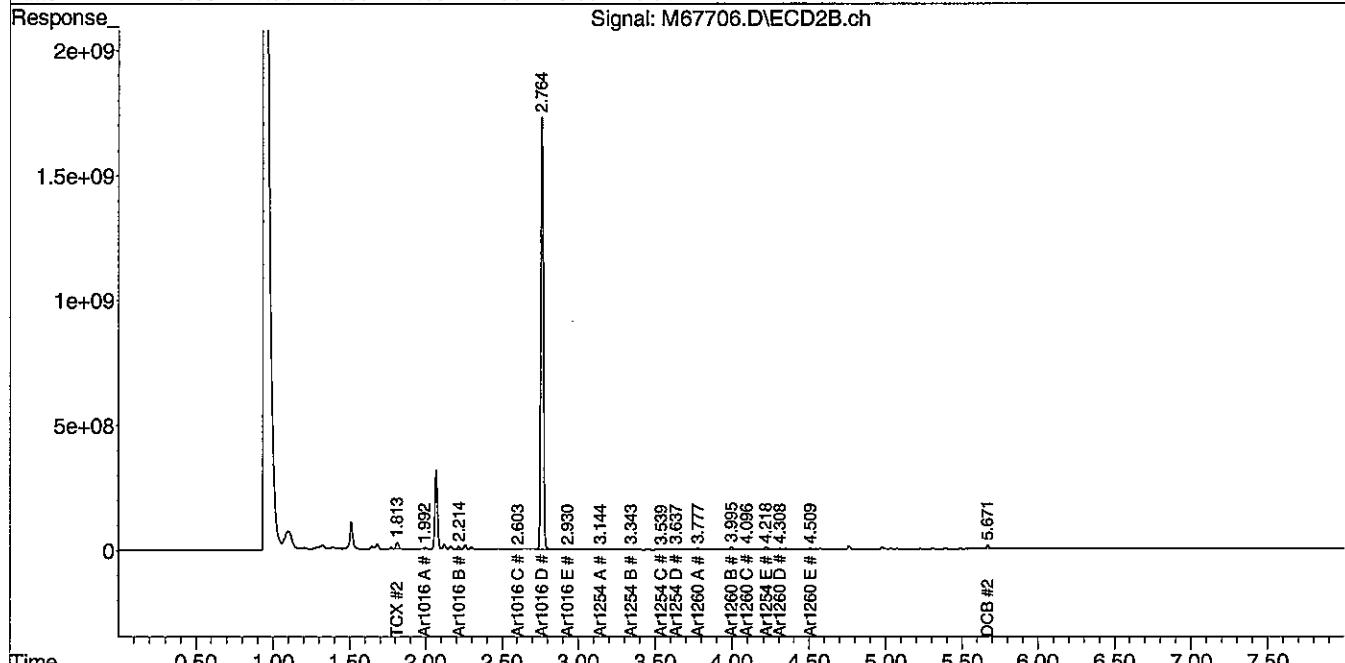
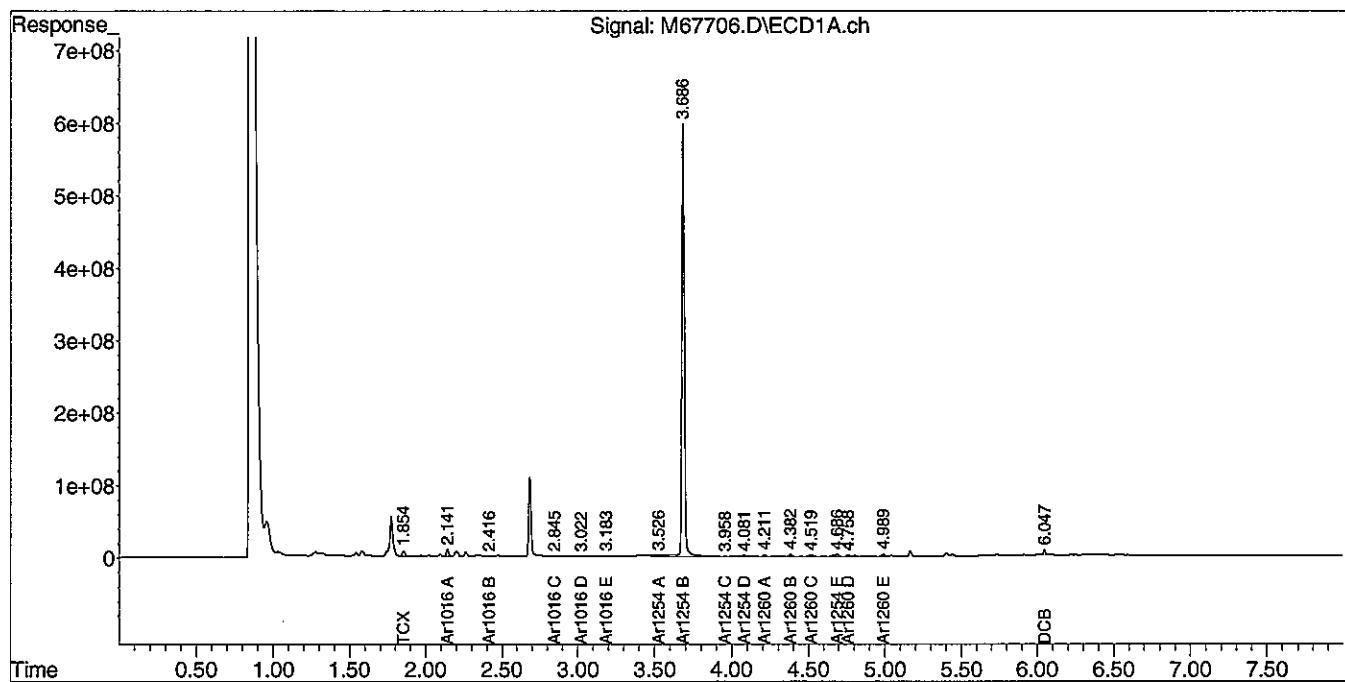
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67706.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 1:22 am
 Operator : JK
 Sample : 74886-15,1:2
 Misc : SOIL
 ALS Vial : 50 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:51 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-229 (0-2')

Lab Sample ID: 74886-16
Matrix: Solid
Percent Solid: 67
Dilution Factor: 1420
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	46900	U
PCB-1221	46900	U
PCB-1232	46900	U
PCB-1242	46900	U
PCB-1248	46900	U
PCB-1254	46900	U
PCB-1260	46900	577000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-16,1:1000

Column ID: 0.25 mm

Data File: M67707.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1414.8

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	571783	577041	0.9	

Column to be used to flag RPD values greater than QC limit of 40%

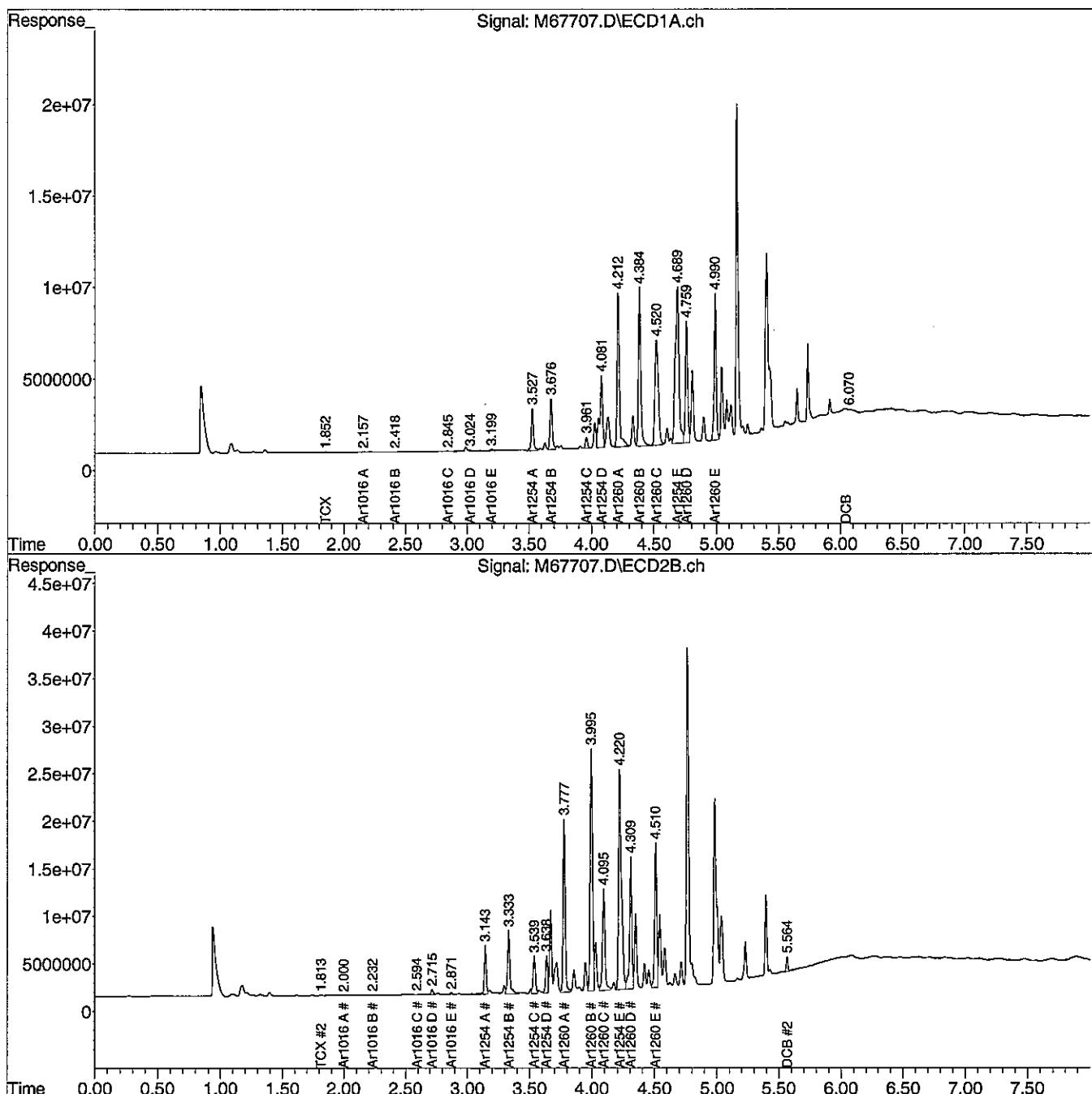
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67707.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 1:32 am
 Operator : JK
 Sample : 74886-16,1:1000
 Misc : SOIL
 ALS Vial : 51 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:53 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-229 (2-4")

Lab Sample ID:	74886-17
Matrix:	Solid
Percent Solid:	87
Dilution Factor:	565
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	18600	U
PCB-1221	18600	U
PCB-1232	18600	U
PCB-1242	18600	U
PCB-1248	18600	U
PCB-1254	18600	U
PCB-1260	18600	192000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-17,1:500

Column ID: 0.25 mm

Data File: M67708.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 565.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	191908	191124	0.4	

Column to be used to flag RPD values greater than QC limit of 40%

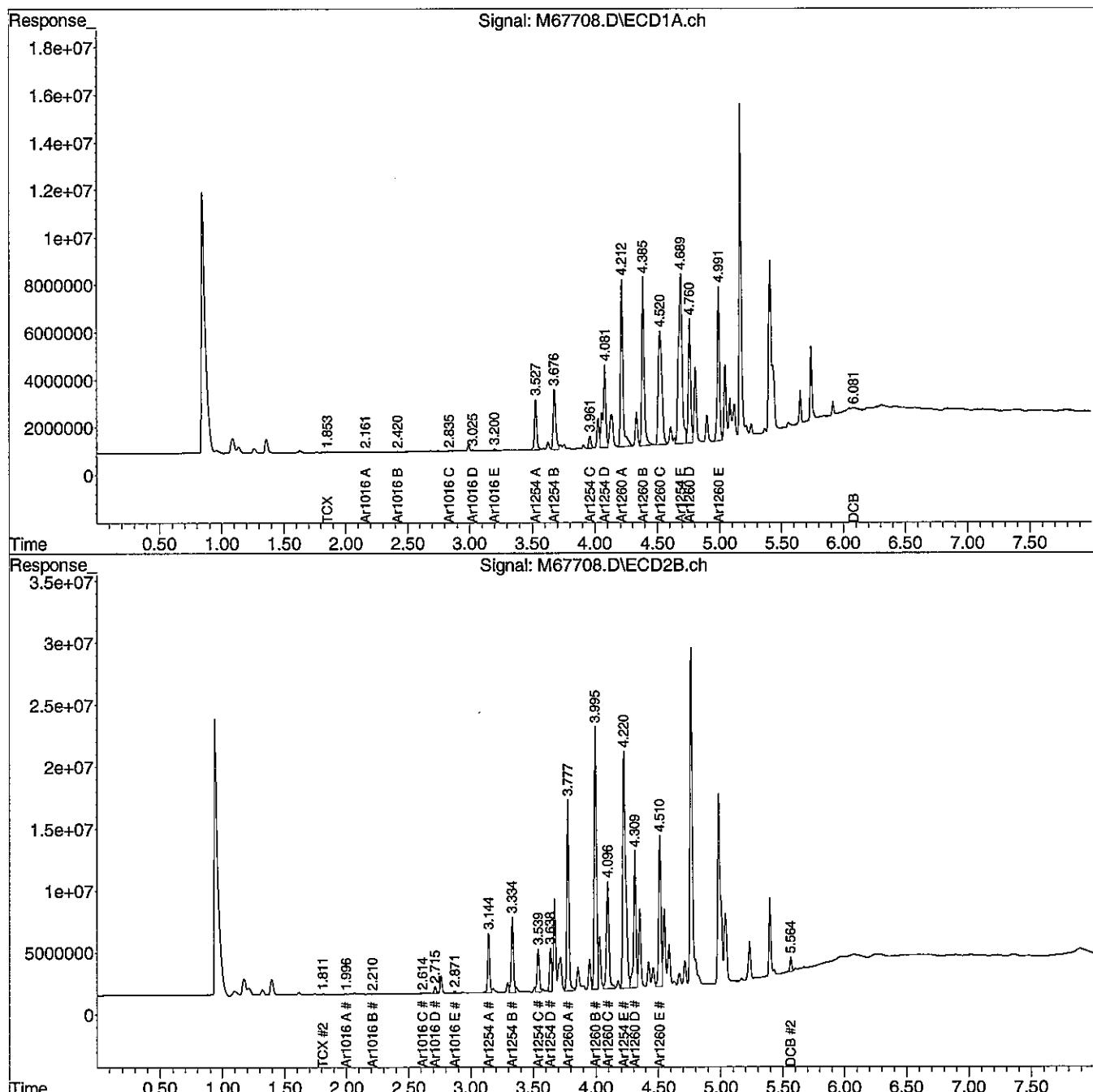
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67708.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 1:42 am
 Operator : JK
 Sample : 74886-17,1:500
 Misc : SOIL
 ALS Vial : 52 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:55 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-229 (4-6')

Lab Sample ID:	74886-18
Matrix:	Solid
Percent Solid:	90
Dilution Factor:	106
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/26/13
Analysis Date:	02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	3500	U
PCB-1221	3500	U
PCB-1232	3500	U
PCB-1242	3500	U
PCB-1248	3500	U
PCB-1254	3500	U
PCB-1260	3500	50600

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB Report

Authorized signature

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-18,1:100

Column ID: 0.25 mm

Data File: M67631.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 106.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	50569	47392	6.5	

Column to be used to flag RPD values greater than QC limit of 40%

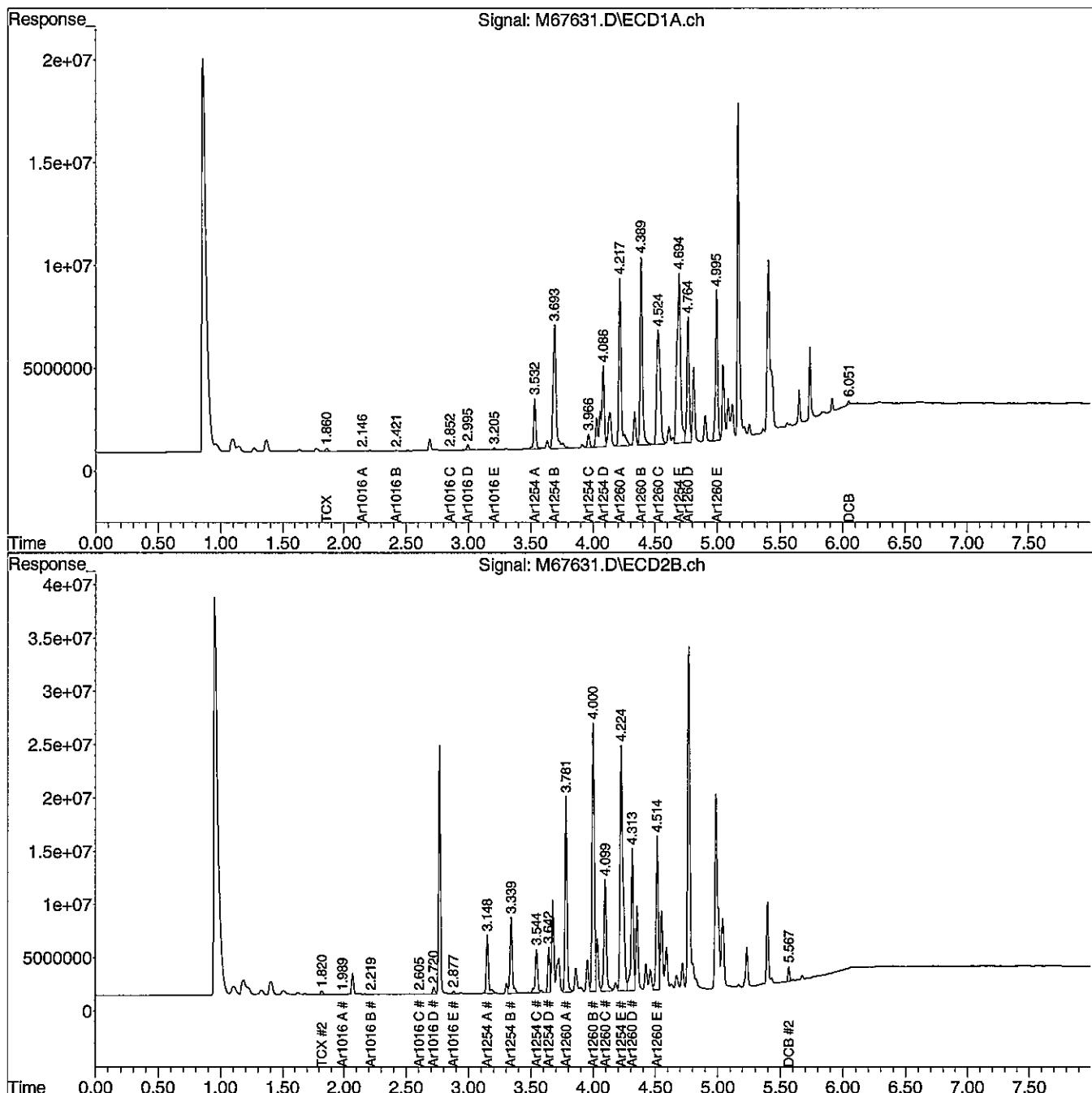
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67631.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:54 am
 Operator : JK
 Sample : 74886-18,1:100
 Misc : SOIL
 ALS Vial : 9 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:51:59 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY

Project Number: 12-3259.1

Field Sample ID: SE-SB-229 (6-8')

Lab Sample ID: 74886-19
Matrix: Solid
Percent Solid: 87
Dilution Factor: 215
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/26/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	7100	U
PCB-1221	7100	U
PCB-1232	7100	U
PCB-1242	7100	U
PCB-1248	7100	U
PCB-1254	7100	U
PCB-1260	7100	149000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB Report

Authorized signature



PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-19,1:200

Column ID: 0.25 mm

Data File: M67632.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 215.3

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	148790	139320	6.6	

Column to be used to flag RPD values greater than QC limit of 40%

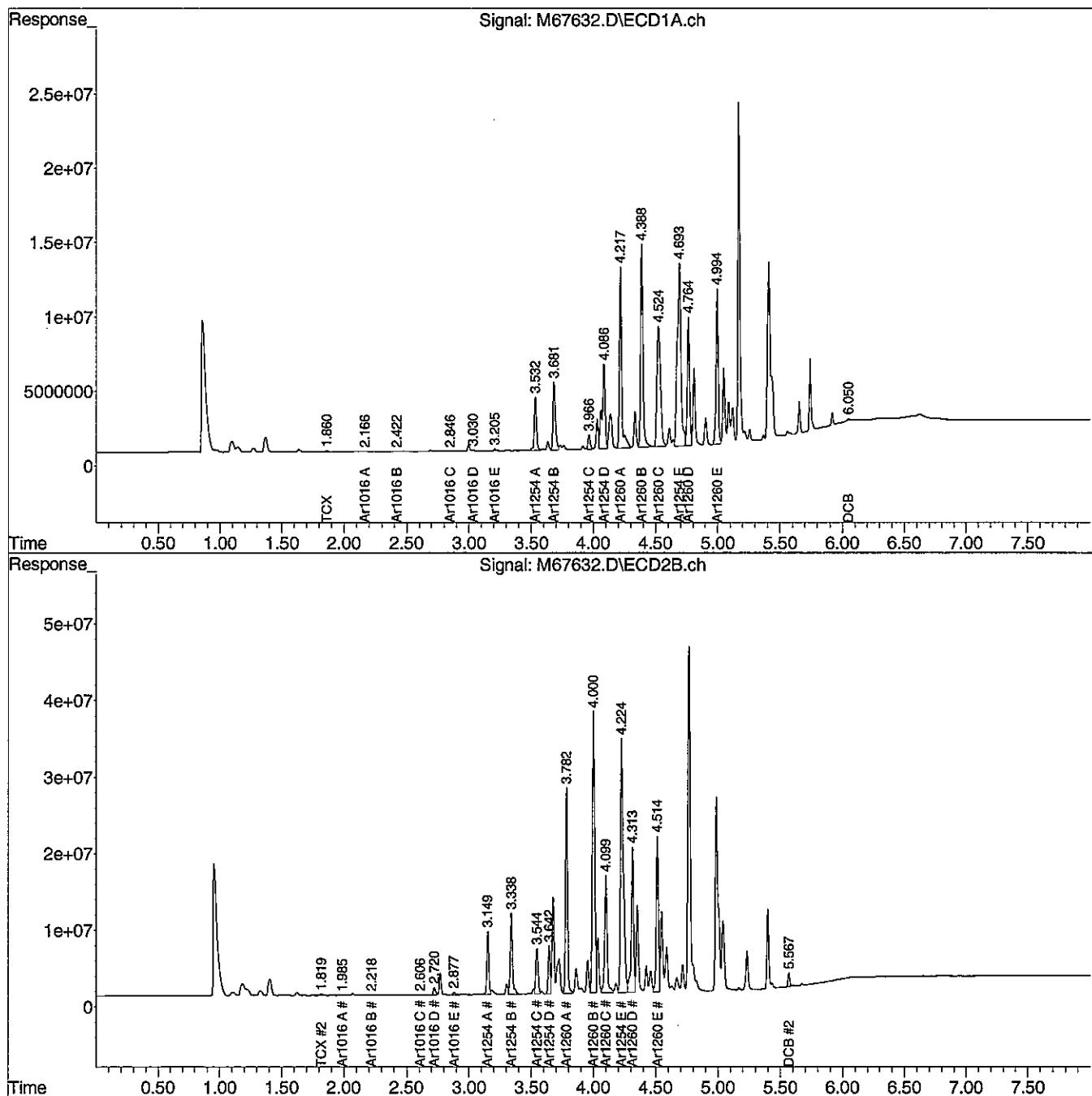
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67632.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 12:04 pm
 Operator : JK
 Sample : 74886-19,1:200
 Misc : SOIL
 ALS Vial : 10 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:52:01 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
 Project Number: 12-3259.1
 Field Sample ID: SE-SB-229 (8-10')

Lab Sample ID: 74886-20
 Matrix: Solid
 Percent Solid: 88
 Dilution Factor: 56
 Collection Date: 02/20/13
 Lab Receipt Date: 02/22/13
 Extraction Date: 02/26/13
 Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	1850	U
PCB-1221	1850	U
PCB-1232	1850	U
PCB-1242	1850	U
PCB-1248	1850	U
PCB-1254	1850	U
PCB-1260	1850	26300

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74886

GC Column #1: STX-CLPesticides I

Sample: 74886-20,1:50

Column ID: 0.25 mm

Data File: M67633.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 55.9

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	26309	25303	3.9	

Column to be used to flag RPD values greater than QC limit of 40%

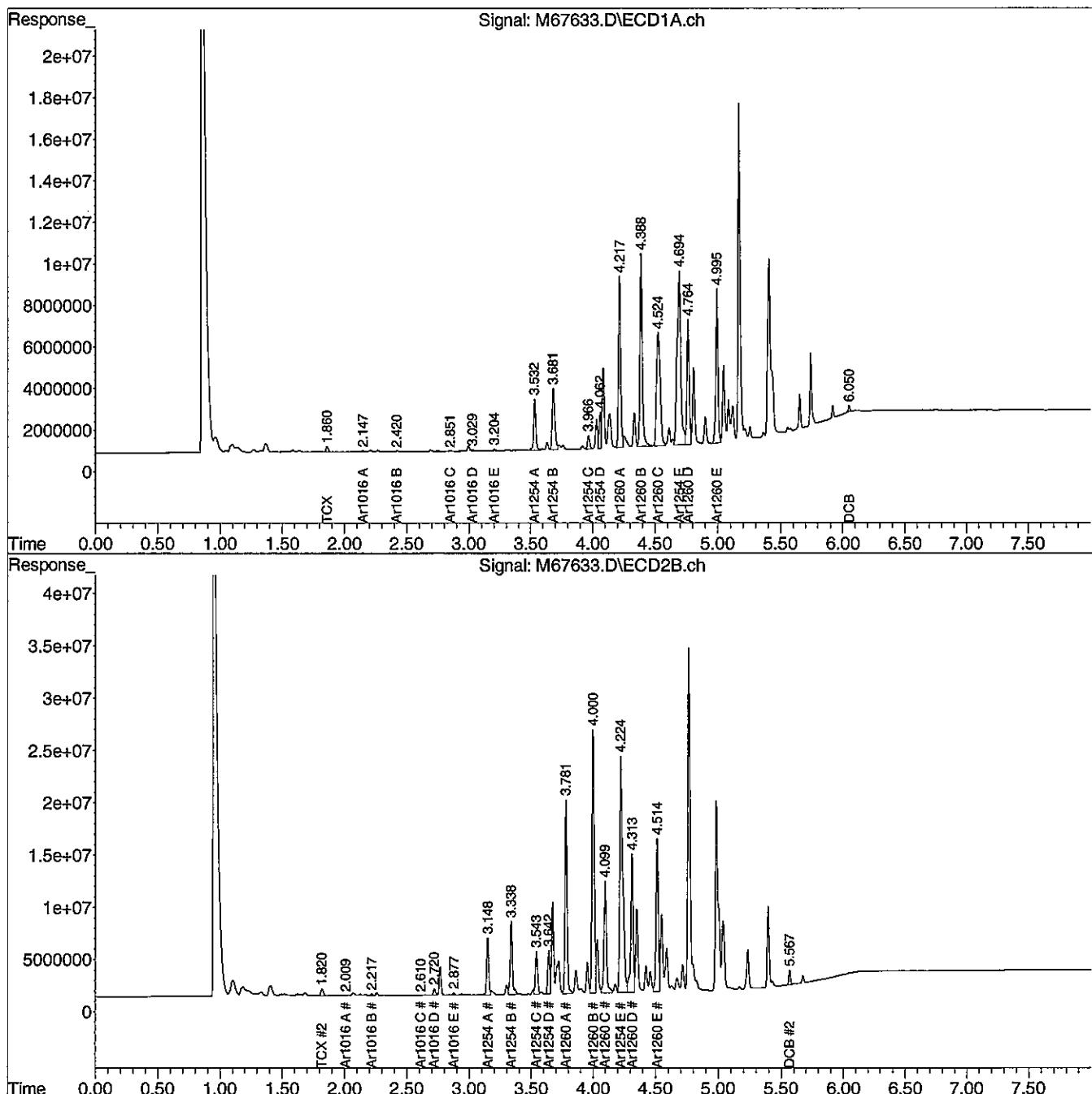
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67633.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 12:14 pm
 Operator : JK
 Sample : 74886-20,1:50
 Misc : SOIL
 ALS Vial : 11 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:52:03 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



PCB
QC FORMS

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 Lewiston ME 04240

February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022613PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/26/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	77	%	
Decachlorobiphenyl	79	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A.
 Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

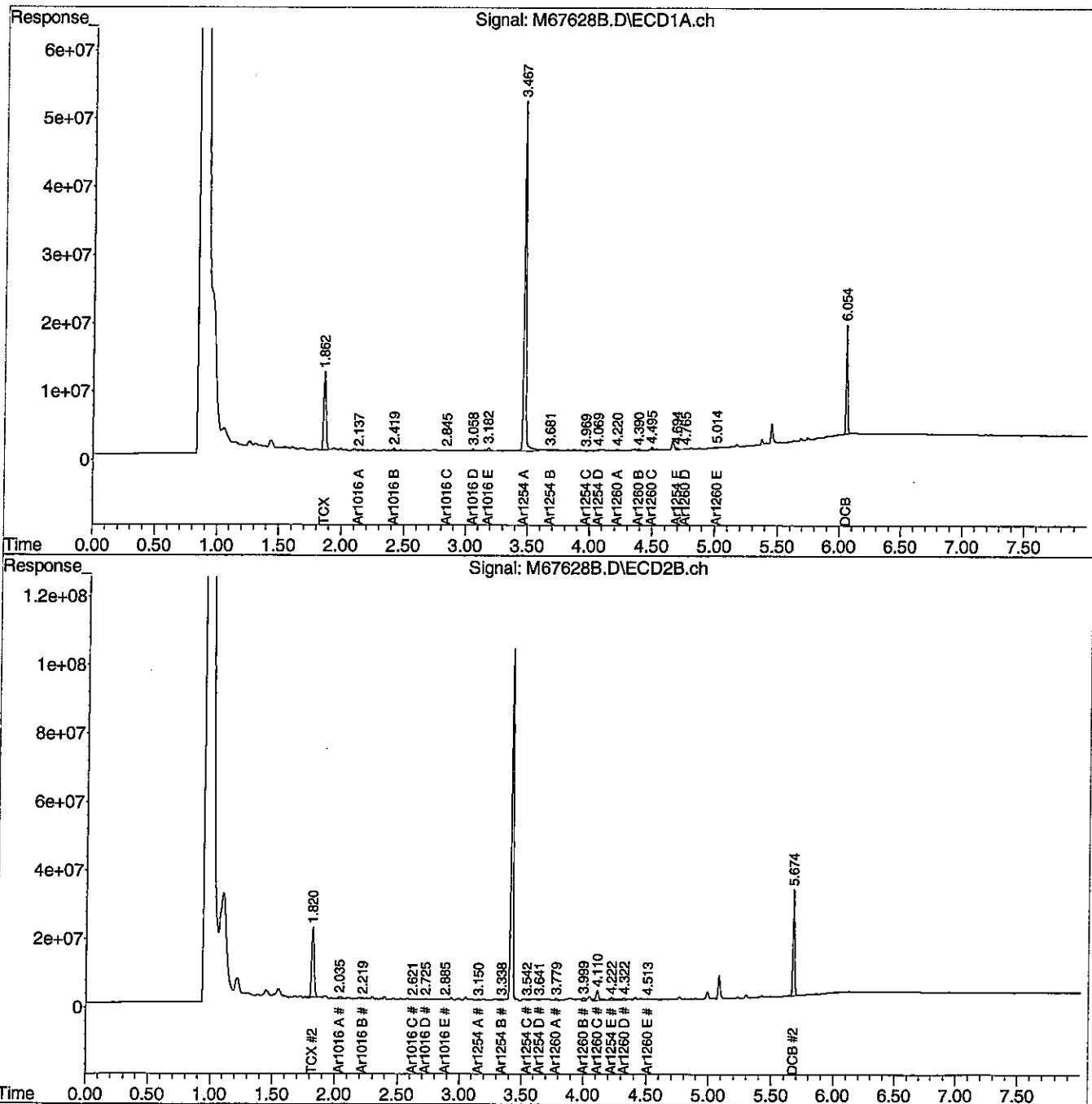
PCB Report

Authorized signature

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67628B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 11:24 am
 Operator : JK
 Sample : B022613PSOX
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:51:53 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:11 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022213PSOX2
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	82	%
Decachlorobiphenyl	78	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

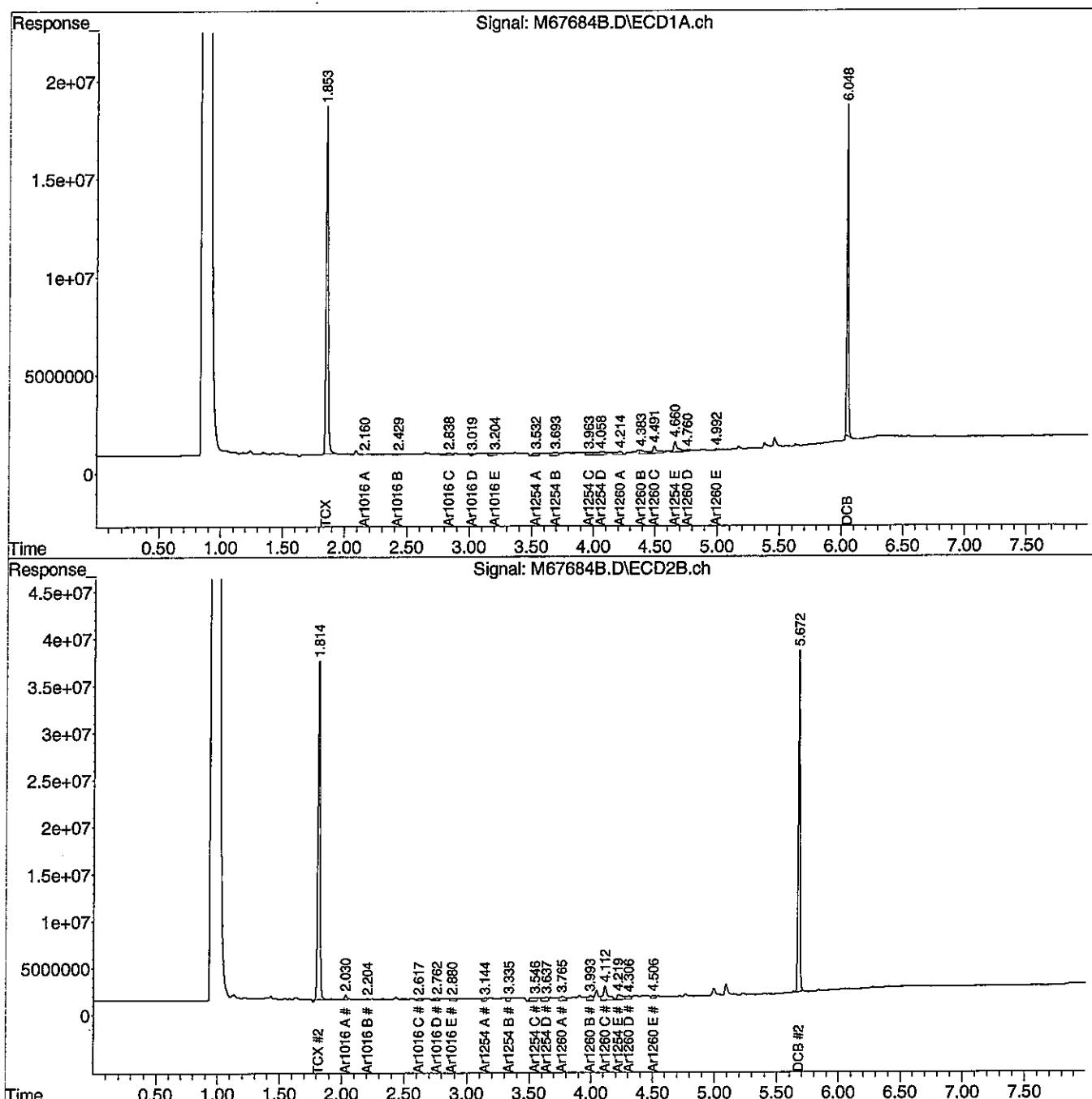
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67684B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 9:40 pm
 Operator : JK
 Sample : B022213PSOX2,,A/C
 Misc : SOIL
 ALS Vial : 28 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:06 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:39 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022513PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/22/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	63	%
Decachlorobiphenyl	78	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

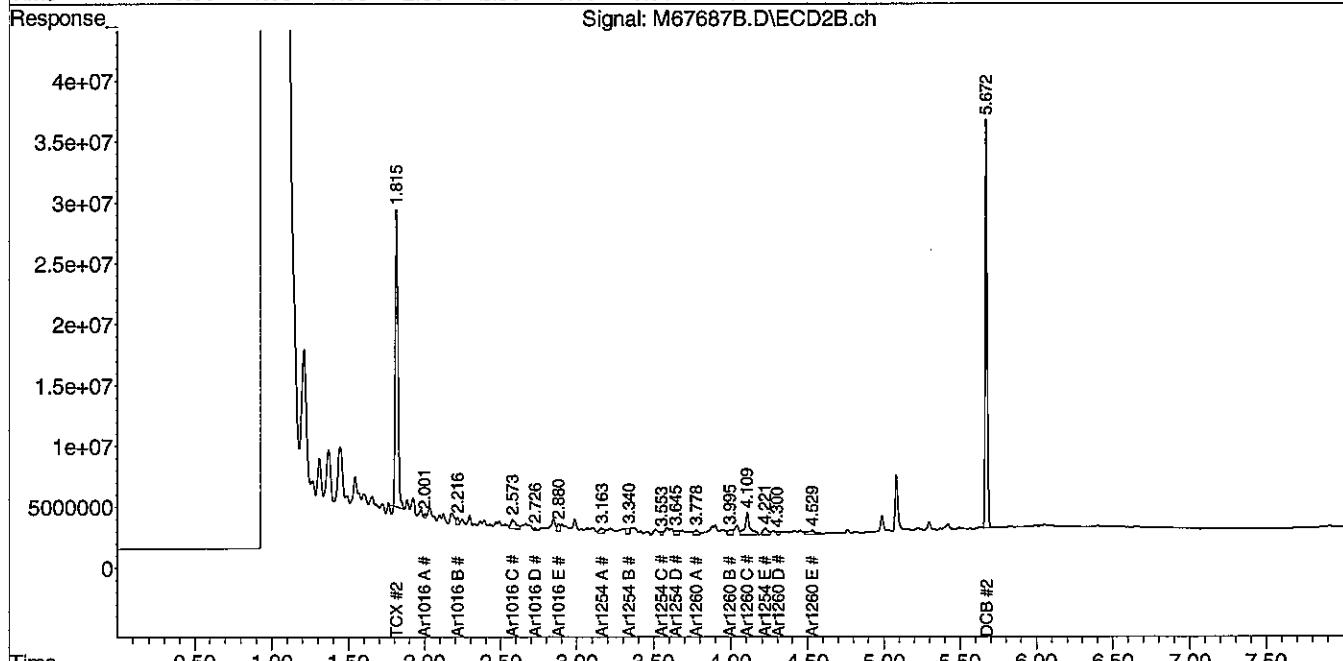
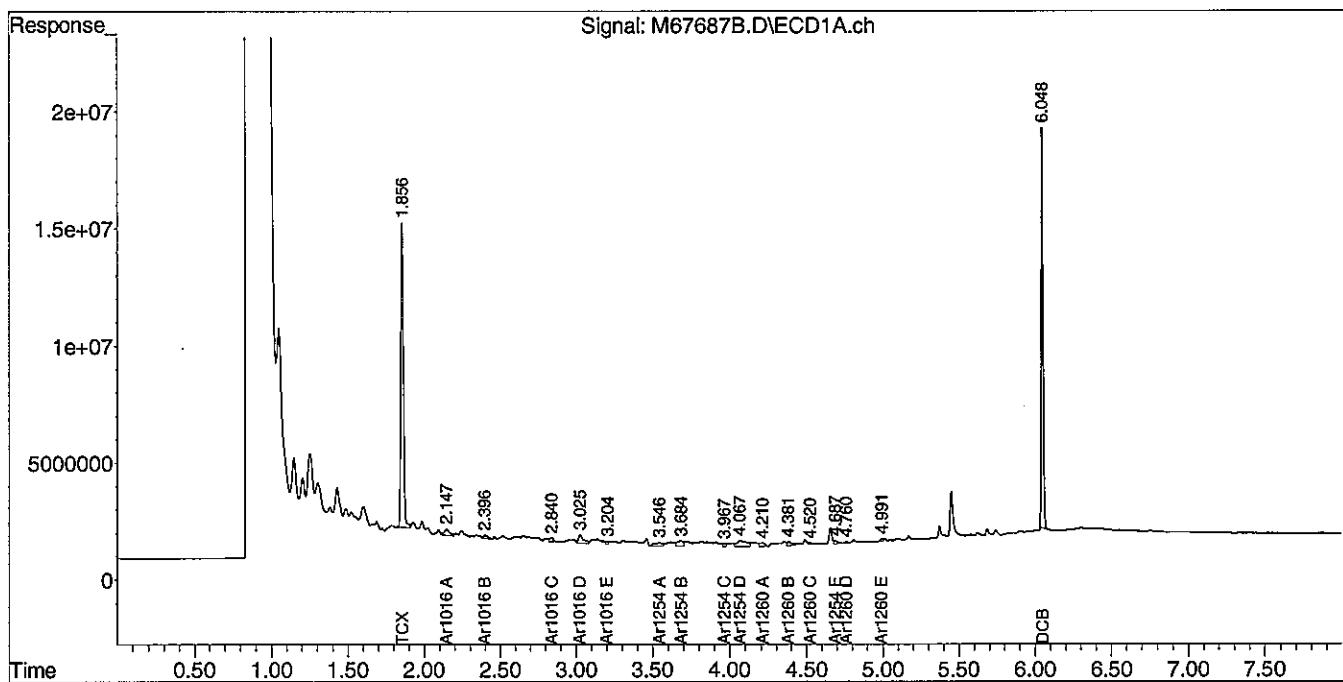
Authorized signature



Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67687B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 10:10 pm
 Operator : JK
 Sample : B022513PSOX
 Misc : SOIL
 ALS Vial : 31 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 09:23:13 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022813PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/28/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	88	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

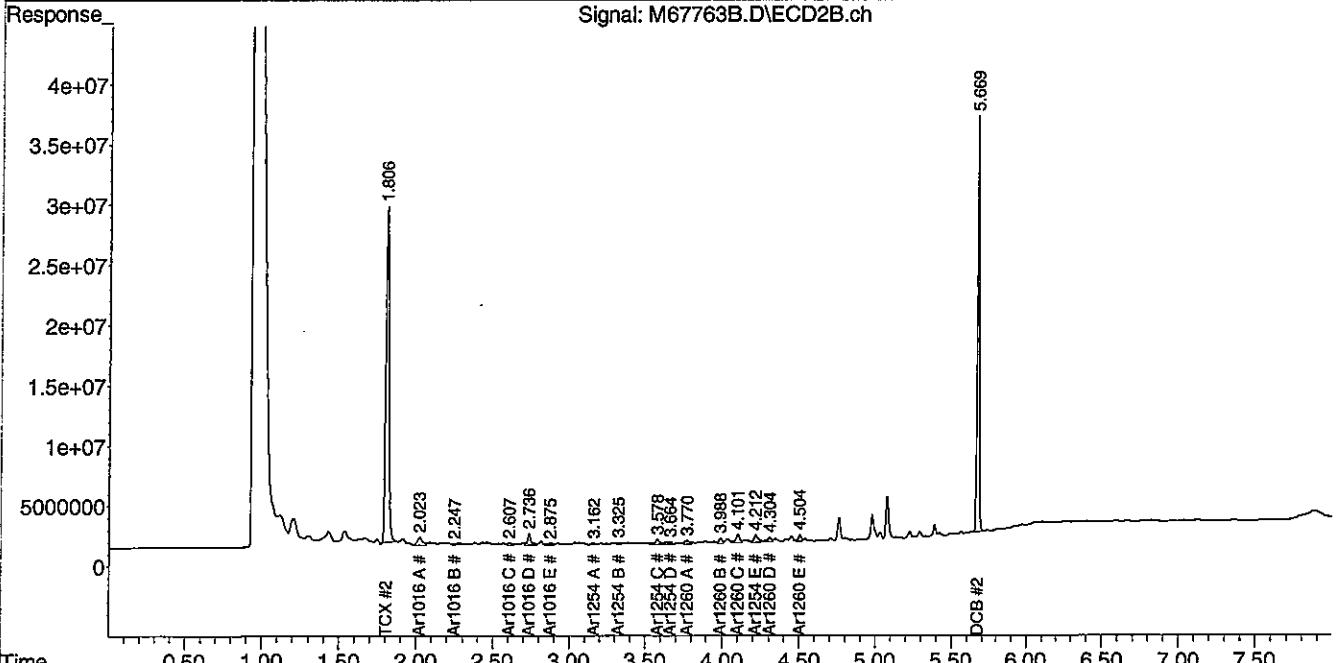
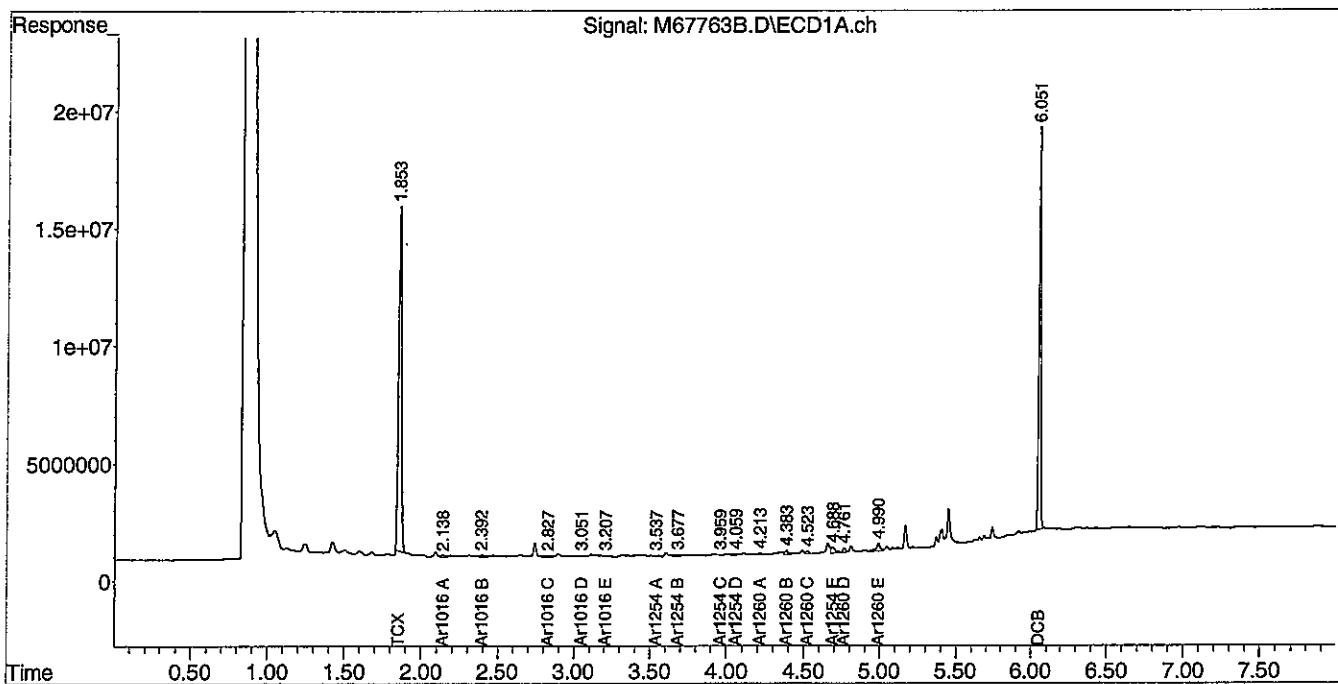
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\023013-M\
 Data File : M67763B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 11:58 am
 Operator : JK
 Sample : B022813PSOX,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 15:34:30 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:11 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74886

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74886

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Lower Limit Upper Limit

SMC #1 = TCX

40 130

SMC #2 = DCB

40 130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74886

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022213PSOX2,,A/C

Spike: L022213PSOX2,,A/C

Spike duplicate: LD022213PSOX2,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	174	87		179	89		3.0	
PCB 1260	200	200	60	130	30	0	178	89		177	88		0.6	
PCB 1016 #2	200	200	65	140	30	0	173	86		172	86		0.3	
PCB 1260 #2	200	200	60	130	30	0	195	97		197	99		1.1	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022513PSOX

Spike: L022513PSOX

Spike duplicate: LD022513PSOX

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	161	81		164	82		2.1	
PCB 1260	200	200	60	130	30	0	170	85		173	87		2.1	
PCB 1016 #2	200	200	65	140	30	0	165	83		165	82		0.3	
PCB 1260 #2	200	200	60	130	30	0	182	91		190	95		4.6	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74886

Column ID: 0.25 mm

Non-spiked sample: 74886-1,1:10.,A/C

GC Column #2: STX-CLPesticides II

Spike: 74886-1,MS,1:10.,A/C

Column ID: 0.25 mm

Spike duplicate: 74886-1,MSD,1:10.,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC #	RESULT (ug/kg)	% REC #	RPD #
PCB 1016	216	224	65	140	30	0	249	115	230	103	7.8
PCB 1260	216	224	60	130	30	2345	6109	1742	*	4488	958 *
PCB 1016 #2	216	224	65	140	30	0	242	112	269	120	10.5
PCB 1260 #2	216	224	60	130	30	2354	5924	1653	*	4396	913 *

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022613PSOX

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: L022613PSOX

Spike duplicate: LD022613PSOX

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	181	90		195	97		7.5	
PCB 1260	200	200	60	130	30	0	182	91		187	94		2.9	
PCB 1016 #2	200	200	65	140	30	0	167	83		179	90		7.3	
PCB 1260 #2	200	200	60	130	30	0	192	96		189	94		1.7	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form

Project Name: MAINE ENERGY		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-3906		(603) 436-5111 (603) 430-2151 Fax		For Analytics Use Only <i>Samples were shipped or hand-delivered 4/0</i>				
Project #: 12-3257.1	Preservation Code:	Circle and/or Write Required Analysis Followed by Preservation Code Please <input checked="" type="checkbox"/> In preservation code here				1) Shipped in good condition: <input checked="" type="checkbox"/> or N 2) Temperature (°C): _____ 3) Received in good condition: <input checked="" type="checkbox"/> or N 4) pH checked by: <i>EG 02/22/13</i> 5) Labels checked by: <i>EG 02/22/13</i>				
Company: SULLIVAN CRESSY	Preservation Key: A = HCl B = 4 C C = Unpres D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other					Matrix Key: C = Concrete WP = Wine WW = Wastewater SW = Surface Water E = Extract				
Report to: 640 MAIN ST. Lewiston, ME 04240						GW = Groundwater DW = Drinking Water S = Soil / Sludge O = Oil X = Other				
Address: 640 MAIN ST. Lewiston, ME 04240										
Phone: (207) 795-6009										
Quote #:										
PO# (if required): 12-3259.1										
Sample Identification	Sample Date	Sample Time					Matrix	No. of Containers checked	Analytics Sample #	
SE-13-234 (3-11)	3-20	1138	N				S	1	74886-1	
SE-13-234 (4-6)	2-20	1131	N				S	1	-2	
SE-13-234 (6-8)	2-20	1133	I				S	1	-3	
SE-13-234 (8-10)	2-20	1154	I				S	1	-4	
SE-13-230 (5-3)	2-20	1156	N				S	1	-5	
SE-13-230 (2-4)	2-20	1203	N				S	1	-6	
SE-13-231 (5-3)	2-20	1215	N				S	1	-7	
SE-13-231 (3-4)	2-20	1220	N				S	1	-8	
SE-13-231 (4-5)	2-20	1224	N				S	1	-9	
SE-13-231 (4-8)	2-20	1235	N				S	1	-10	
Comments, Additional Analyses, or Special Instructions:										** List requested metals here *Fee may apply
Email Results to: <i>XCEENY@summitenergy.com</i>	Turnaround Time (TAT)		Report Type:		State Standard:					
<input type="checkbox"/> 1 Day*	<input type="checkbox"/> 2 Days*	<input type="checkbox"/> 3 Days*	<input checked="" type="checkbox"/> 4 Days*	<input type="checkbox"/> Standard (4-10 business days)	<input type="checkbox"/> MCP*	<input type="checkbox"/> Level II*	<input type="checkbox"/> NH	<input type="checkbox"/> MA	<input type="checkbox"/> ME	
<input type="checkbox"/> 5 Days					<input type="checkbox"/> CTRCP*	<input type="checkbox"/> Level III*	<input type="checkbox"/> CT	<input type="checkbox"/> RI	<input type="checkbox"/> EDD Required: Y* <input checked="" type="checkbox"/> (eg. S-1 or GW-1)	
					<input type="checkbox"/> DOD*	<input type="checkbox"/> Level IV*	<input type="checkbox"/> Standard	<input type="checkbox"/> Other	<input type="checkbox"/> Type: _____	
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.										
Sampler Name (Print): <i>John Cressy</i>	Date: <i>2-21</i>	Time: <i>8:00</i>	Received By: <i>General Counsel</i>							
Relinquished By Sampler: <i>John Cressy</i>	Date: <i>2/21/13</i>	Time: <i>8:50</i>	Received By: <i>Jeffrey</i>							
Relinquished By: <i>John Cressy</i>	Date: <i>2/21/13</i>	Time: <i>8:50</i>	Received By: <i>Jeffrey</i>							

Chain Of Custody Form



environmental
laboratory LLC

Project Name: MARIE ENERG Project#: 12-3359.1		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9306		(603) 436-5111 (603) 430-2151 Fax	
For Analytics Use Only					
Samples were: 1) Shipped or hand-delivered 40 2) Temperature (°C): 0 3) Received in good condition: Y br N 4) pH checked by: Fee3 02/22/13 5) Labels checked by: Fee3 02/22/13					
Matrix Key: C = Concrete WP = Wipe WW = Wastewater SW = Surface Water E = Extract					
Preservation Key: A = HCl B = 4°C C = Unpreserved D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other					
Metals: CRAs PPI3 TAL23 Other** VPH: Full or Ranges only TEPH EPH: Full or Ranges only ME4125 TPB: 8015 (Diesel Range) 8100M ME4125 TPH: 8015 (Gas Range) ME4125 PCB: 8082 608 Solvent V.O. Pesticides: 8081 608 VOC: 8270 625 PAH only SIM SVOC: 8260 524.2 624 Field Filtered Y N					
Sample Identification	Sample Date	Sample Time	Please circle and/or write required analysis followed by preservation code here	Please circle and/or write required analysis followed by preservation code here	No. of containers checked
SE-SB-235 (0-2)	2-20	1250	X	X	1
SE-SB-235 (0-4)	2-20	1254	X	X	1
SE-SB-235 (4-6)	2-20	1300	N	X	1
SE-SB-235 (6-8)	2-20	1302	N	X	1
SE-SB-235 (8-10)	2-20	1311	N	X	1
SE-SB-239 (0-2)	2-20	1325	N	X	1
SE-SB-239 (2-4)	2-20	1329	N	X	1
SE-SB-239 (4-6)	2-20	1330	N	X	1
SE-SB-239 (6-8)	2-20	1335	N	X	1
SE-SB-239 (8-10)	2-20	1340	N	X	1
					-20
Comments, Additional Analyses, or Special Instructions: * Cap notches coc but label states "SE-SB-234 (2-4)" with sample date of 12/20/13 → COC + cap correct per client Fee3					
** List requested metals here Project Requirements: *Fee may apply					
Report Type: <input type="checkbox"/> MCP* <input type="checkbox"/> Level II* <input type="checkbox"/> MA <input type="checkbox"/> ME <input type="checkbox"/> CTRCP* <input type="checkbox"/> Level III* <input type="checkbox"/> CT <input type="checkbox"/> RI <input type="checkbox"/> DOD* <input type="checkbox"/> Level IV* <input type="checkbox"/> RI <input checked="" type="checkbox"/> Standard <input type="checkbox"/> Other: Fee3 Type: _____					
Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 4 Days* <input type="checkbox"/> 3 Days* <input checked="" type="checkbox"/> Standard (3-10 business days)					
*Fee may apply; lab approval required					
Relinquished By Sampler: John Cresset Relinquished By: John Cresset Relinquished By:					
Date: 3-21 Time: 9:00 Received By: John Cresset Date: 3/22/13 Time: 8:50 Received By: John Cresset Date:					



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 74886
CLIENT: Summit
PROJECT: Maine Energy

COOLER NUMBER: 255
NUMBER OF COOLERS: 1

A: PRELIMINARY EXAMINATION:

1. Cooler received by(initials): JB

DATE COOLER RECEIVED/OPENED: 2/22/13

2. Circle one: Hand delivered
(If so, skip 3)

Shipped

3. Did cooler come with a shipping slip?

Y

N

3a. Enter carrier name and airbill number here:
504730

4. Were custody seals on the outside of cooler?

How many & where: _____ Seal Date: _____ Seal Name: _____

Y

N

5. Did the custody seals arrive unbroken and intact upon arrival?

Y

N/A

6. COC#:

7. Were Custody papers filled out properly (ink,signed, legible, project information etc)?

Y

N

8. Were custody papers sealed in a plastic bag?

Y

N

9. Did you sign the COC in the appropriate place?

Y

N

10. Was enough ice used to chill the cooler?

Y N Temp. of cooler: 4°C

B. Log-In: Date samples were logged in:

2/22/13

By: JB

11. Were all bottles sealed in separate plastic bags?

Y

N

12. Did all bottles arrive unbroken and were labels in good condition?

Y

N

13. Were all bottle labels complete(ID,Date,time,etc.)

Y

N

14. Did all bottle labels agree with custody papers?

Y

N

see coc

15. Were the correct containers used for the tests indicated:

Y

N

16. Were samples received at the correct pH?

Y

N/A

17. Was sufficient amount of sample sent for the tests indicated?

Y

N

18. Were all samples submitted within holding time?

Y

N

19. Were all containers used within AEL's expiration date?**

Y

N

20. Were VOA samples absent of greater than pea-sized bubbles?

Y

N/A

(Note:Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

*If NO, List Sample ID's, Lab #s:

When bubbles are present in VOA samples they are labelled from smallest (or no bubbles) to largest. Lab to analyze VOA samples with no bubbles or smallest bubbles first

20. Laboratory labeling verified by (initials): CCJ

Date: 02/22/13

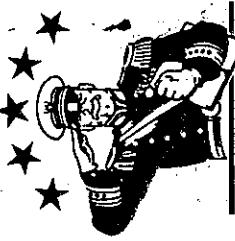
**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.

Insured Delivery

P.O. Box 1072
Portland, ME 04104

385 Main Street
So. Portland, ME 04106
1-800-698-5035
767-6004.

FAX: 767-7159
E-mail: misraelson@generalcourier.com



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R			
O			
M			

Summit Environmental

640 Main St
Lewiston, Me

Analytics Lab

195 Commerce Way
Parsonsfield, N.H.

BILLING INFORMATION

QUANTITY	DESCRIPTION	PRICE
65	856	Time 856 2/22/13

Signature

Jaclyn

March 1, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74887
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples
PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
Chromatograms
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:
No exceptions.

PCBs by EPA Method 8082:
No results were reported below the quantitation limit.

All samples except 74887-13, 74887-14 and 74887-21 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument.

The MS/MSD analyzed on sample 74887-18 did not meet acceptance criteria for PCB 1260 recoveries due to the parent sample having concentrations of PCB 1260 that exceeded the calibration range of the instrument. The laboratory control samples (L022253PSOX2/LD022253PSOX2) were in control for all recoveries and RPDs. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC



Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
Summit Environmental Consultants Inc.
640 Main Street
Lewiston ME 04240

Report Number: 74887

Revision: Rev. 0

Re: MAINE ENERGY (Project No: 12-3259.1)

Enclosed are the results of the analyses on your sample(s). Samples were received on 22 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

Sample Analysis: The attached pages detail the Client Sample IDs, Lab Sample IDs, and Analyses requested

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

If you have any questions on these results, please do not hesitate to contact us.

Authorized signature



Stephen L. Knollmeyer Lab. Director

Date

3/1/2013

This report shall not be reproduced, except in full, without the written consent of Analytics Environmental Laboratory, LLC.

**CLIENT: Summit Environmental
 Consultants Inc.**

REPORT NUMBER: 74887

REV: Rev. 0

PROJECT: MAINE ENERGY (Project No: 12-3259.1)

<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74887-1	02/20/13	SE-SB-227 (0-2')	EPA 8082 (PCBs only)	
74887-2	02/20/13	SE-SB-227 (2-4')	EPA 8082 (PCBs only)	
74887-3	02/20/13	SE-SB-227 (4-6')	EPA 8082 (PCBs only)	
74887-4	02/20/13	SE-SB-227 (6-8')	EPA 8082 (PCBs only)	
74887-5	02/20/13	SE-SB-228 (0-2')	EPA 8082 (PCBs only)	
74887-6	02/20/13	SE-SB-228 (2-4')	EPA 8082 (PCBs only)	
74887-7	02/20/13	SE-SB-228 (4-6')	EPA 8082 (PCBs only)	
74887-8	02/20/13	SE-SB-228 (6-8')	EPA 8082 (PCBs only)	
74887-9	02/20/13	SE-SB-228 (8-10')	EPA 8082 (PCBs only)	
74887-10	02/20/13	SE-SB-222 (0-2')	EPA 8082 (PCBs only)	
74887-11	02/20/13	SE-SB-222 (2-4')	EPA 8082 (PCBs only)	
74887-12	02/20/13	SE-SB-222 (4-6')	EPA 8082 (PCBs only)	
74887-13	02/20/13	SE-SB-222 (6-8')	EPA 8082 (PCBs only)	
74887-14	02/20/13	SE-SB-222 (8-10')	EPA 8082 (PCBs only)	
74887-15	02/20/13	SE-SB-223 (0-2')	EPA 8082 (PCBs only)	
74887-16	02/20/13	SE-SB-223 (2-4')	EPA 8082 (PCBs only)	
74887-17	02/20/13	SE-SB-223 (4-6')	EPA 8082 (PCBs only)	
74887-18	02/20/13	SE-SB-223 (6-8')	EPA 8082 (PCBs only)	
74887-19	02/20/13	SE-SB-223 (8-10')	EPA 8082 (PCBs only)	
74887-20	02/20/13	SE-SB-218 (0-2')	EPA 8082 (PCBs only)	
74887-21	02/20/13	SE-SB-218 (2-4')	EPA 8082 (PCBs only)	
74887-22	02/20/13	SE-SB-231 (8-10')	EPA 8082 (PCBs only)	

Surrogate Compound Limits

	Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water				
1,4-Difluorobenzene		70-130		EPA 524.2
Bromofluorobenzene		70-130		
1,2-Dichlorobenzene-d4		70-130		
Volatile Organic Compounds				
1,2-Dichloroethane-d4		70-120	70-120	EPA 624/8260B
Toluene-d8		85-120	85-120	
Bromofluorobenzene		75-120	75-120	
Semi-Volatile Organic Compounds				
2-Fluorophenol		20-110	35-105	EPA 625/8270C
d5-Phenol		15-110	40-100	
d5-nitrobenzene		40-110	35-100	
2-Fluorobiphenyl		50-110	45-105	
2,4,6-Tribromophenol		40-110	40-125	
d14-p-terphenyl		50-130	30-125	
PAH's by SIM				
d5-nitrobenzene		21-110	35-110	EPA 8270C
2-Fluorobiphenyl		36-121	45-105	
d14-p-terphenyl		33-141	30-125	
Pesticides and PCBs				
2,4,5,6-Tetrachloro-m-xylene (TCX)		46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)		40-135	40-130	
Herbicides				
Dichloroacetic acid (DCAA)		30-150	30-150	
Gasoline Range Organics/TPH Gasoline				
Trifluorotoluene TFT (FID)		60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)		60-140	60-140	
Trifluorotoluene TFT (PID)		60-140	60-140	
Bromofluorobenzene (BFB) (PID)		60-140	60-140	
Diesel Range Organics/TPH Diesel				
m-terphenyl		60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons				
2,5-Dibromotoluene (PID)		70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)		70-130	70-130	
Extricable Petroleum Hydrocarbons				
1-chloro-octadecane (aliphatic)		40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)		40-140	40-140	
2-Fluorobiphenyl (Fractionation)		40-140	40-140	
2-Bromonaphthalene (fractionation)		40-140	40-140	

**PCB
DATA SUMMARIES**

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-227 (0-2")

Lab Sample ID: 74887-1
Matrix: Solid
Percent Solid: 81
Dilution Factor: 12
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	396	U
PCB-1221	396	U
PCB-1232	396	U
PCB-1242	396	U
PCB-1248	396	U
PCB-1254	396	U
PCB-1260	396	2840

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	86	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A.
 Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-1,1;10

Column ID: 0.25 mm

Data File: M67716.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 12.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2749	2841	3.3	

Column to be used to flag RPD values greater than QC limit of 40%

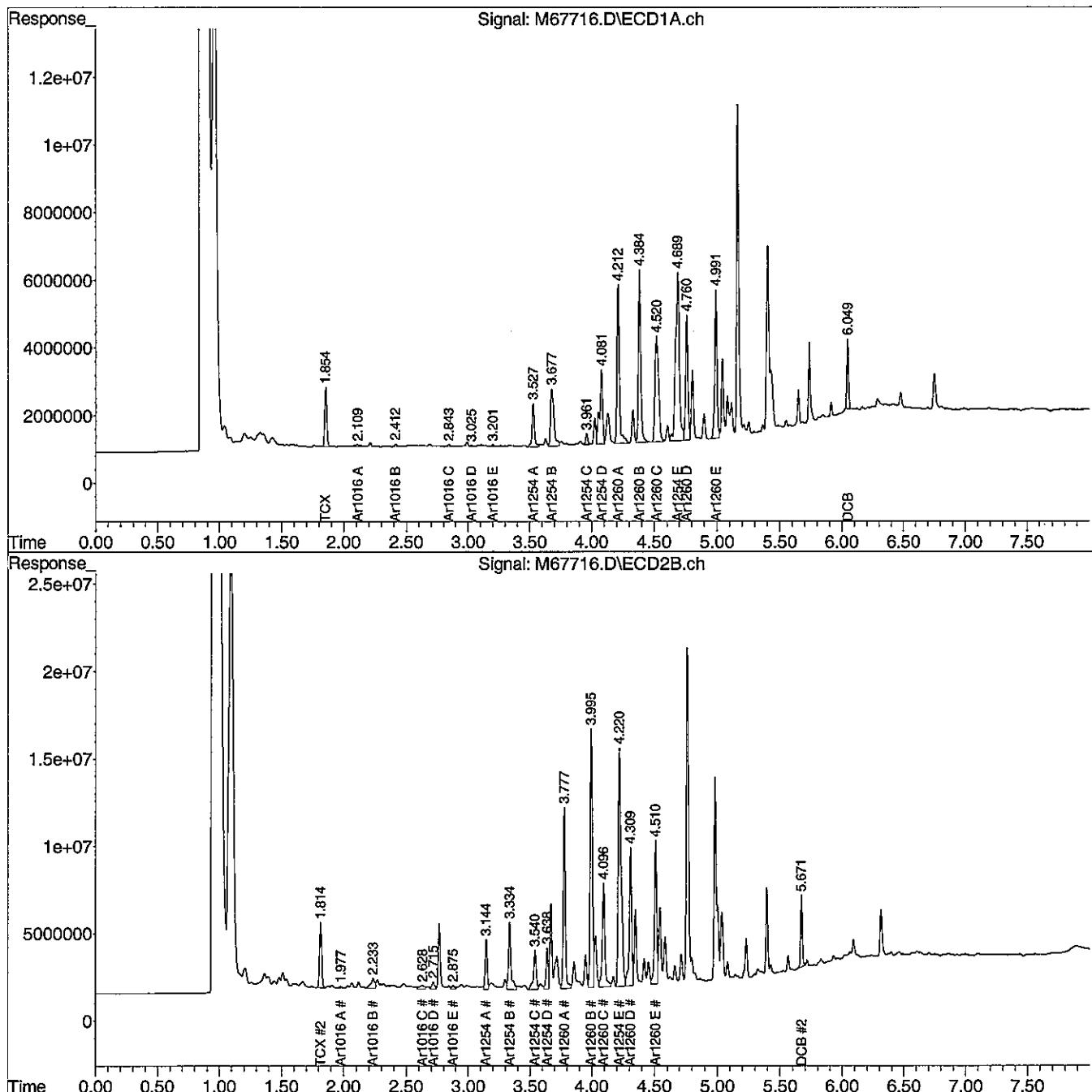
* Values outside QC limits

Comments: _____

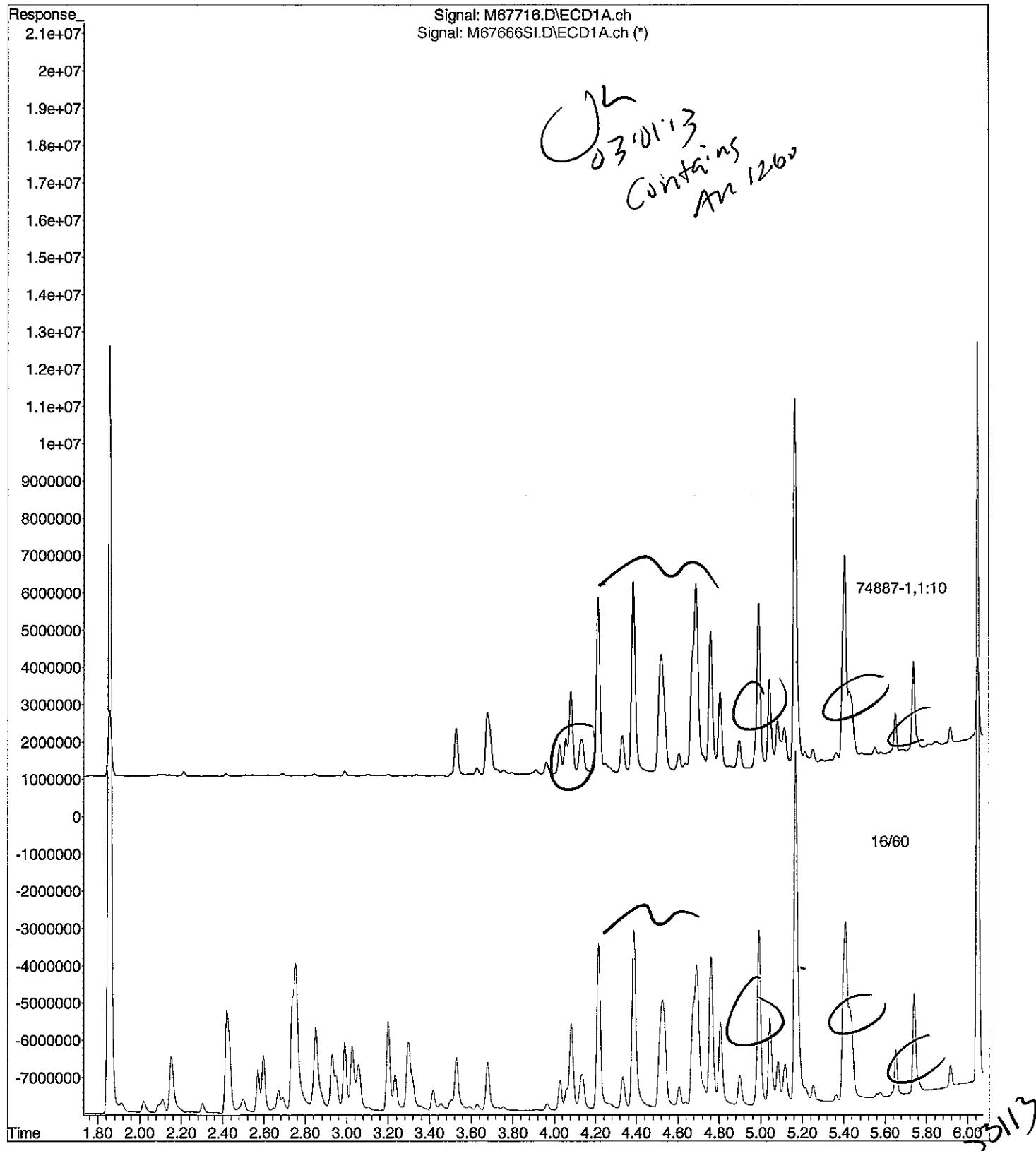
Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67716.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:03 am
 Operator : JK
 Sample : 74887-1,1:10
 Misc : SOIL
 ALS Vial : 56 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:06 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022813-M\M67716.D
Operator : JK
Acquired : 1 Mar 2013 3:03 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74887-1,1:10
Misc Info : SOIL
Vial Number: 56



Mr. John Cressey
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 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-227 (2-4')

Lab Sample ID: 74887-2
Matrix: Solid
Percent Solid: 87
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	1930

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	90	%	
Decachlorobiphenyl	97	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-2,1:10

Column ID: 0.25 mm

Data File: M67717.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1727	1927	11.0	

Column to be used to flag RPD values greater than QC limit of 40%

* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67717.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:13 am
 Operator : JK
 Sample : 74887-2,1:10
 Misc : SOIL
 ALS Vial : 57 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e

Quant Time: Mar 01 10:33:08 2013

Quant Method : C:\msdchem\1\METHODS\PCB022813.M

Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254

QLast Update : Fri Mar 01 08:56:41 2013

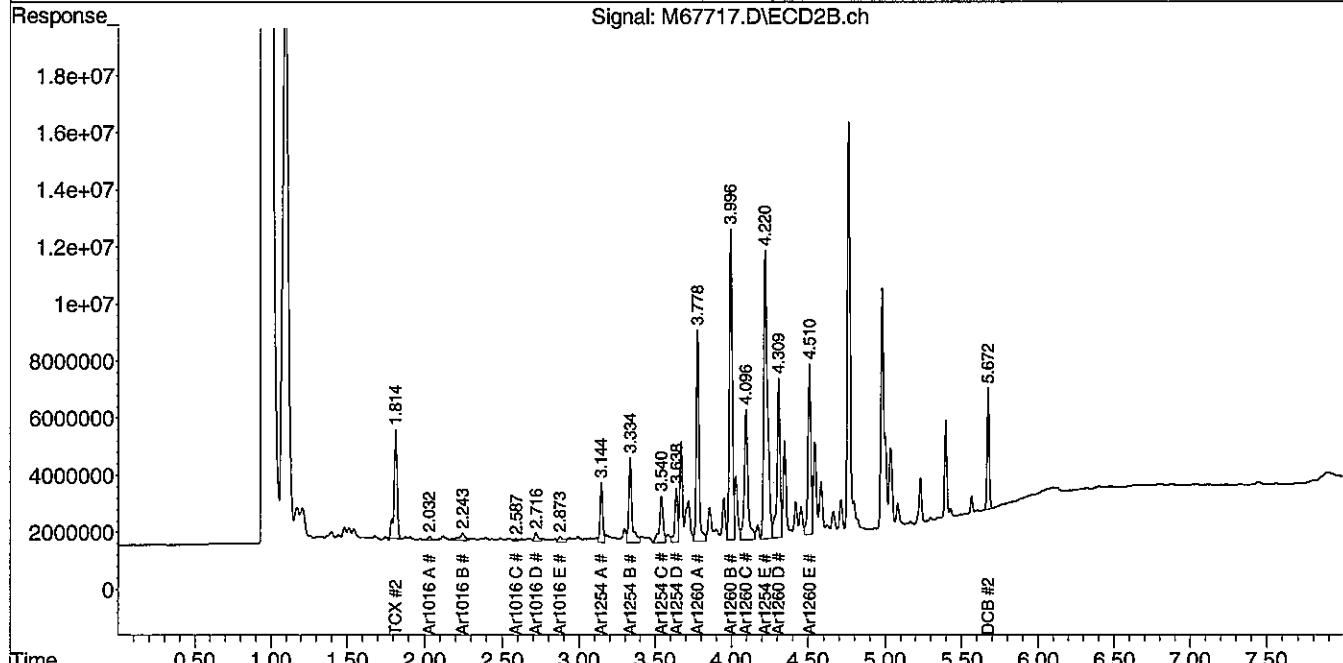
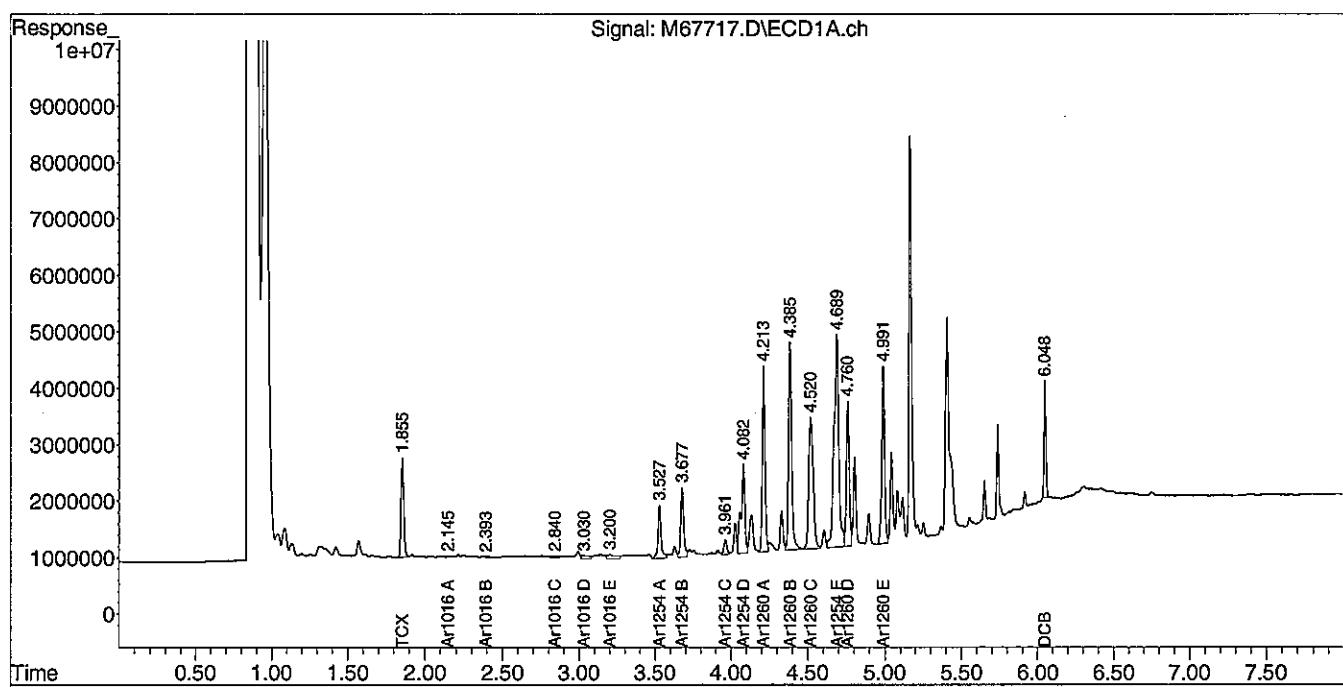
Response via : Initial Calibration

Integrator: ChemStation

Volume Inj. : 2 uL

Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides

Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID	
Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-227 (4-6')

Lab Sample ID: 74887-3
Matrix: Solid
Percent Solid: 87
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	996

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	89	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-3,1:5

Column ID: 0.25 mm

Data File: M67718.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	938	996	6.0	

Column to be used to flag RPD values greater than QC limit of 40%

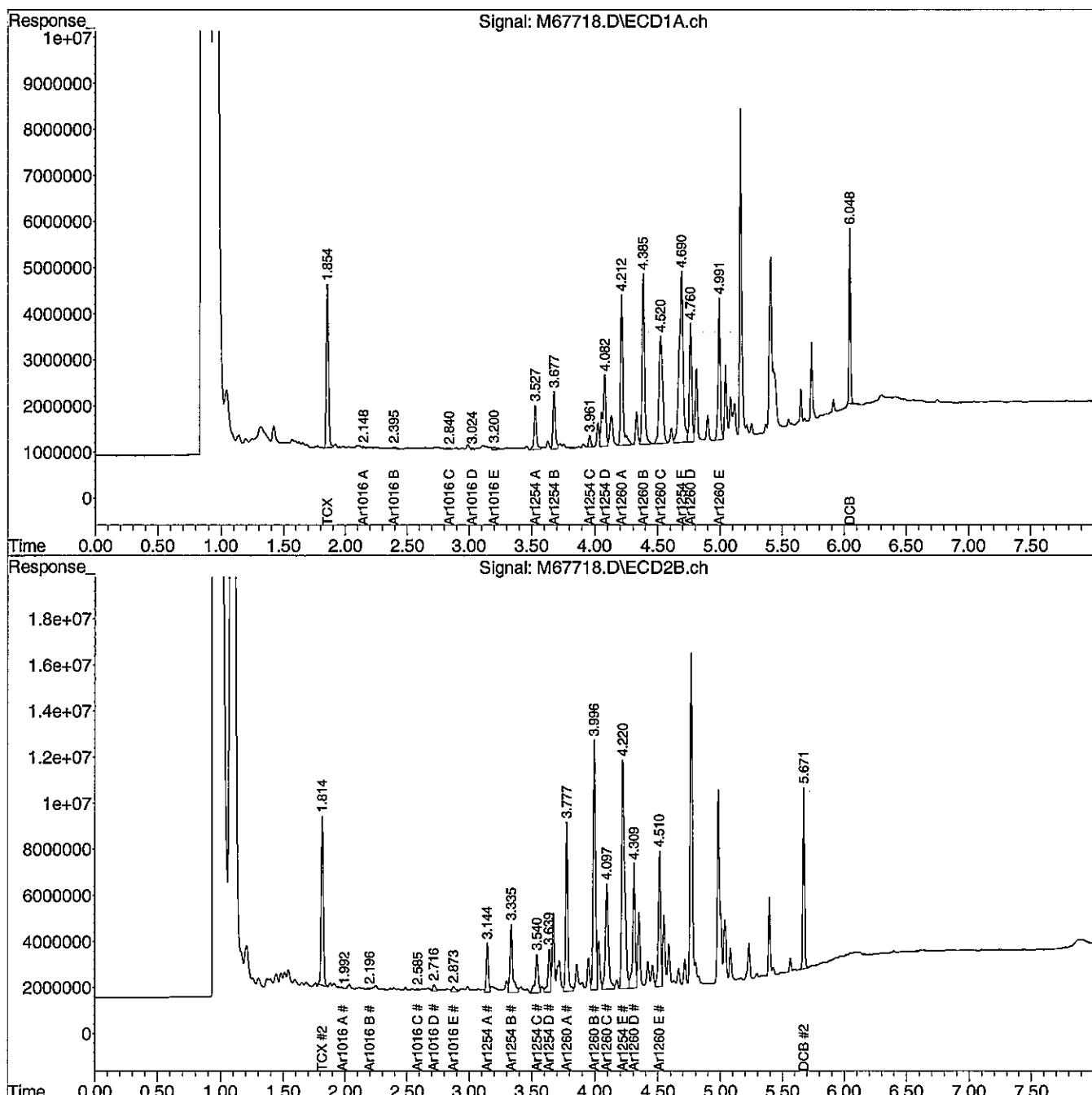
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67718.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:23 am
 Operator : JK
 Sample : 74887-3,1:5
 Misc : SOIL
 ALS Vial : 58 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:10 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 640 Main Street
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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-227 (6-8')

Lab Sample ID: 74887-4
Matrix: Solid
Percent Solid: 90
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	4320

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	86	%
Decachlorobiphenyl	83	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-4,1:10

Column ID: 0.25 mm

Data File: M67719.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	4320	4282	0.9	

Column to be used to flag RPD values greater than QC limit of 40%

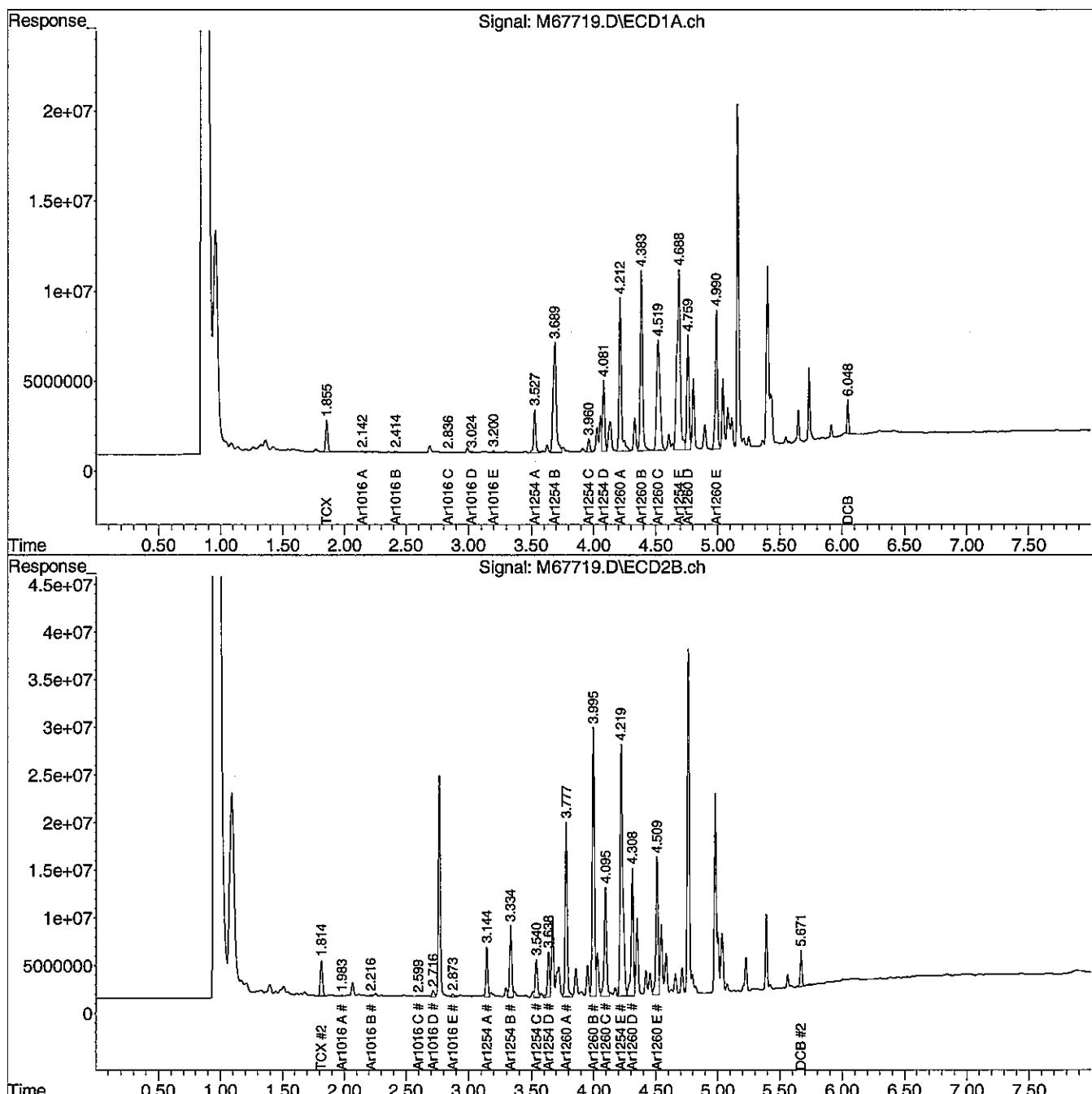
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67719.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:33 am
 Operator : JK
 Sample : 74887-4,1:10
 Misc : SOIL
 ALS Vial : 59 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:12 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-228 (0-2')

Lab Sample ID:	74887-5
Matrix:	Solid
Percent Solid:	84
Dilution Factor:	232
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	7660	U
PCB-1221	7660	U
PCB-1232	7660	U
PCB-1242	7660	U
PCB-1248	7660	U
PCB-1254	7660	U
PCB-1260	7660	86200

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-5,1:200

Column ID: 0.25 mm

Data File: M67720.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 232.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	86248	84887	1.6	

Column to be used to flag RPD values greater than QC limit of 40%

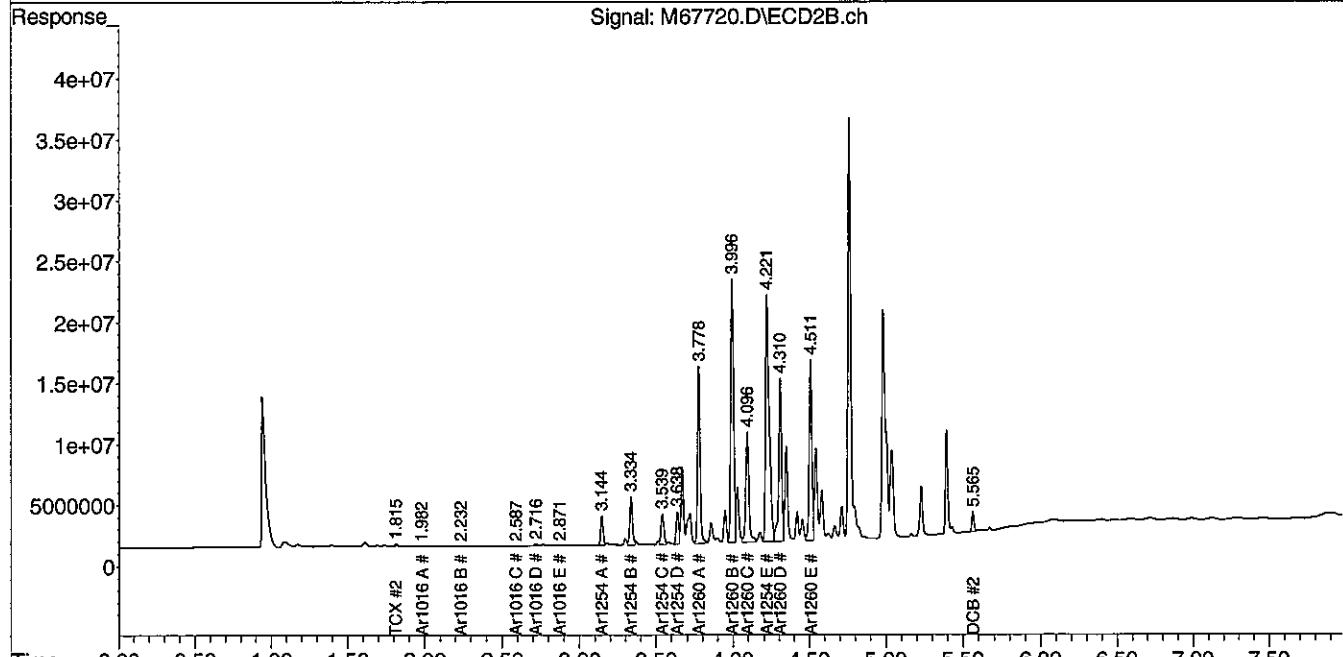
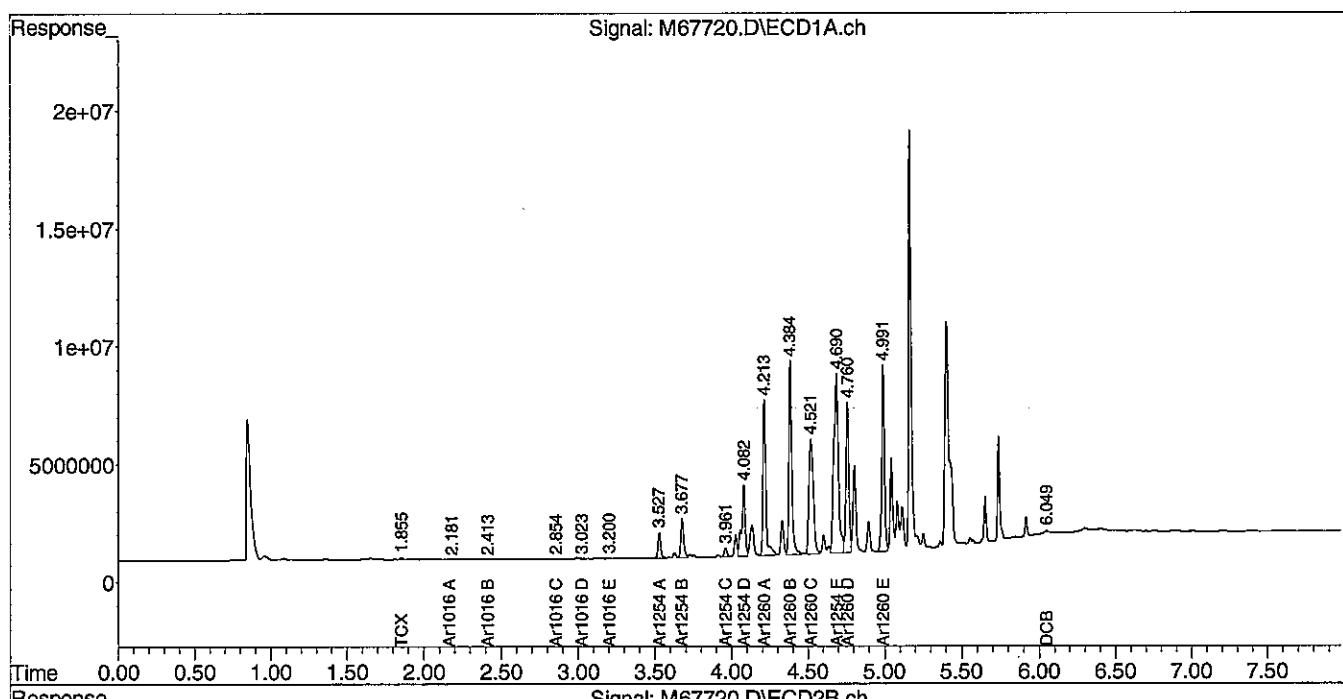
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67720.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:43 am
 Operator : JK
 Sample : 74887-5,1:200
 Misc : SOIL
 ALS Vial : 60 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:14 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-228 (2-4")

Lab Sample ID:	74887-6
Matrix:	Solid
Percent Solid:	87
Dilution Factor:	6
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	1360

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	85	%	
Decachlorobiphenyl	89	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A.
 Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-6,1:5

Column ID: 0.25 mm

Data File: M67721.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.8

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	1264	1356	7.0	

Column to be used to flag RPD values greater than QC limit of 40%

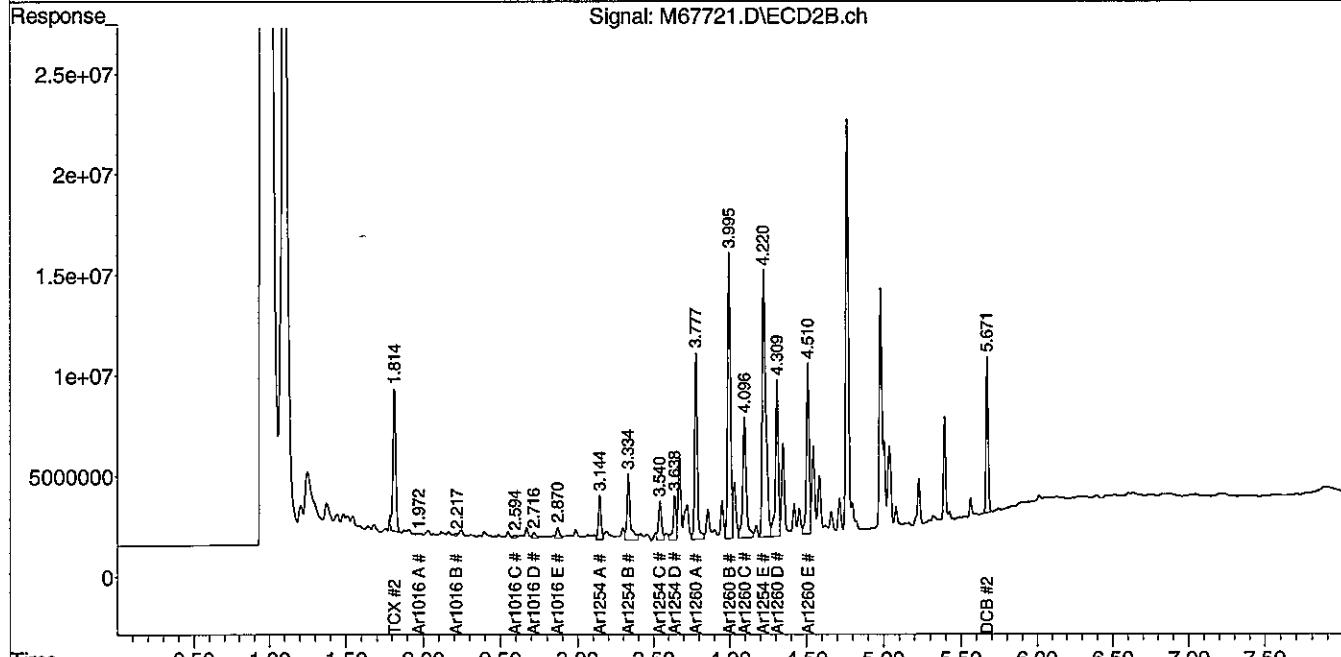
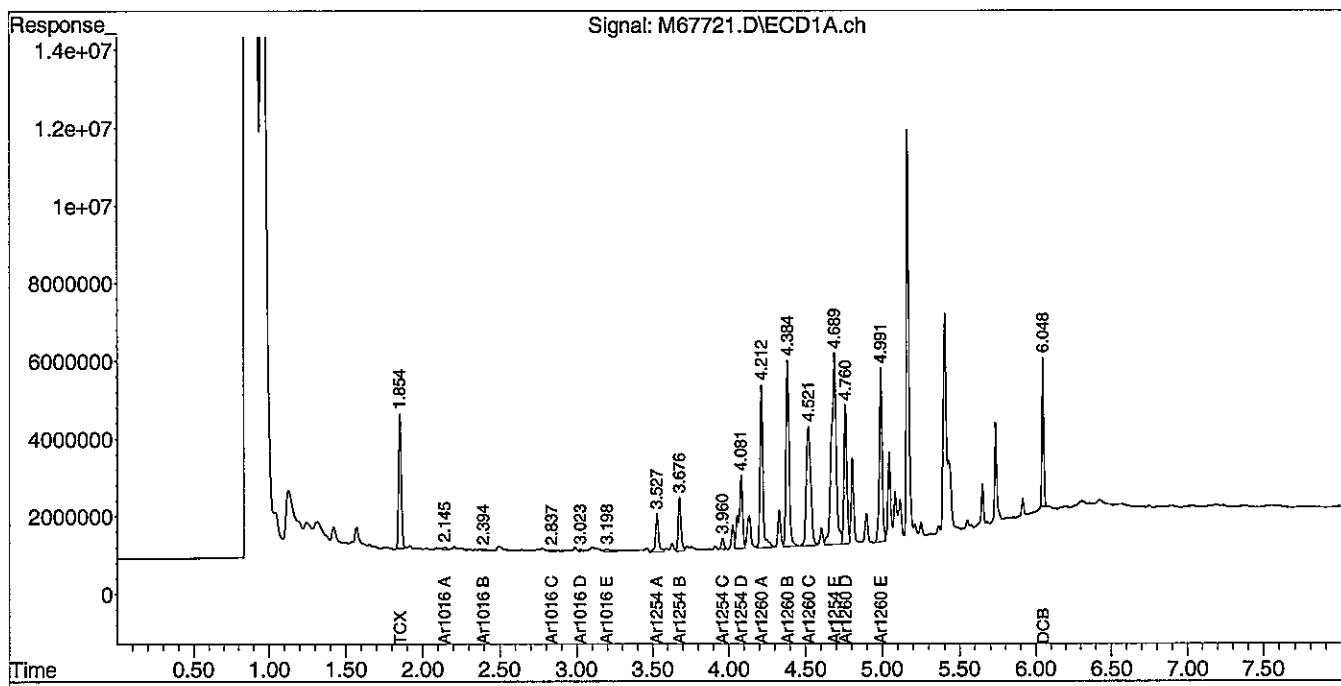
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67721.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 3:53 am
 Operator : JK
 Sample : 74887-6,1:5
 Misc : SOIL
 ALS Vial : 61 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:16 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 640 Main Street
 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-228 (4-6')

Lab Sample ID: 74887-7
Matrix: Solid
Percent Solid: 93
Dilution Factor: 54
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	1780	U
PCB-1221	1780	U
PCB-1232	1780	U
PCB-1242	1780	U
PCB-1248	1780	U
PCB-1254	1780	U
PCB-1260	1780	19500

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-7,1:50

Column ID: 0.25 mm

Data File: M67722.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 53.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	19458	19145	1.6	

Column to be used to flag RPD values greater than QC limit of 40%

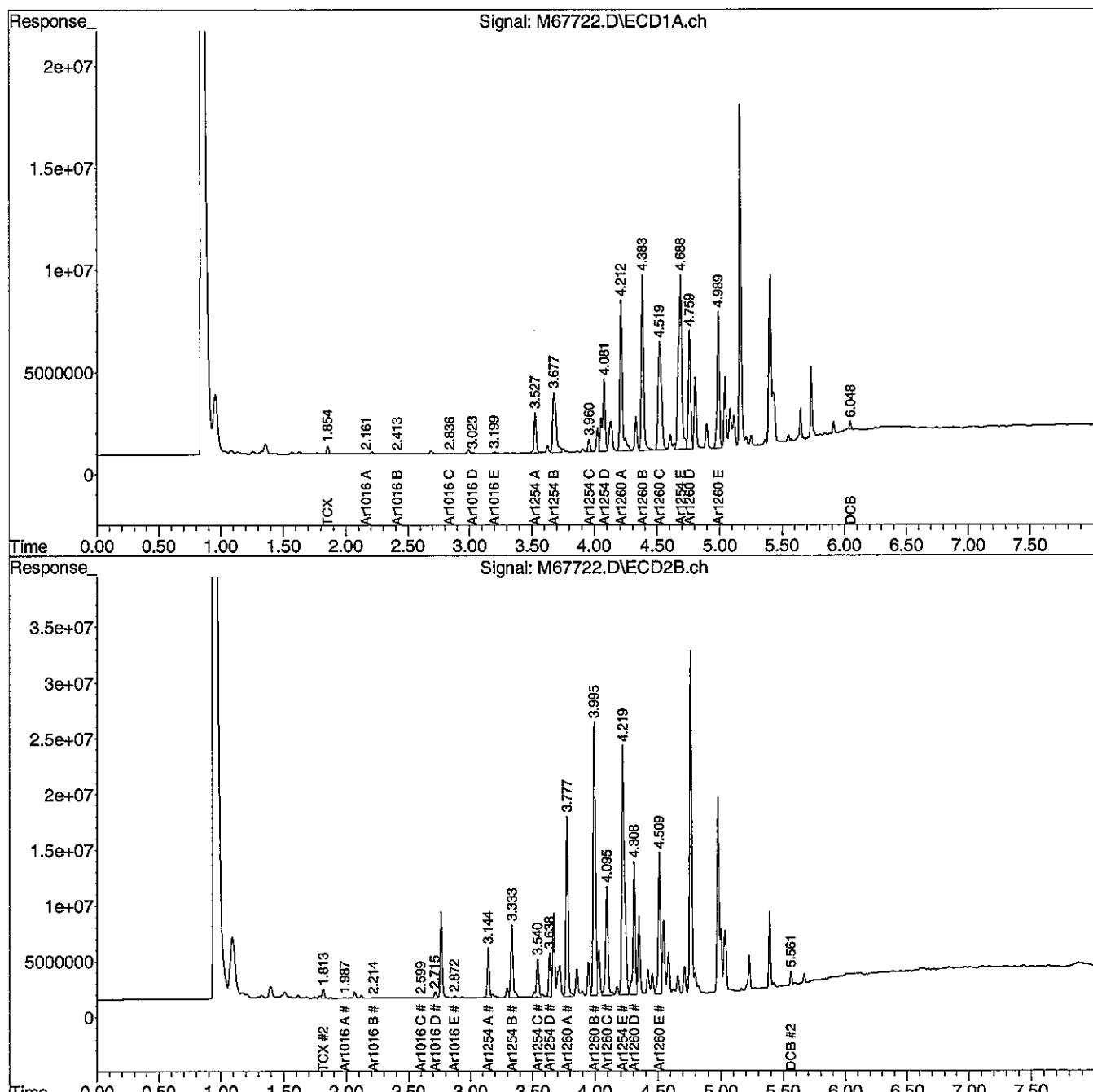
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67722.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:03 am
 Operator : JK
 Sample : 74887-7,1:50
 Misc : SOIL
 ALS Vial : 62 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:18 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



Mr. John Cressey
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 640 Main Street
 Lewiston ME 04240

March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-228 (6-8')

Lab Sample ID: 74887-8
Matrix: Solid
Percent Solid: 91
Dilution Factor: 10
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	330	U
PCB-1221	330	U
PCB-1232	330	U
PCB-1242	330	U
PCB-1248	330	U
PCB-1254	330	U
PCB-1260	330	3980

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	87	%
Decachlorobiphenyl	80	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-8,1:10

Column ID: 0.25 mm

Data File: M67723.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3938	3981	1.1	

Column to be used to flag RPD values greater than QC limit of 40%

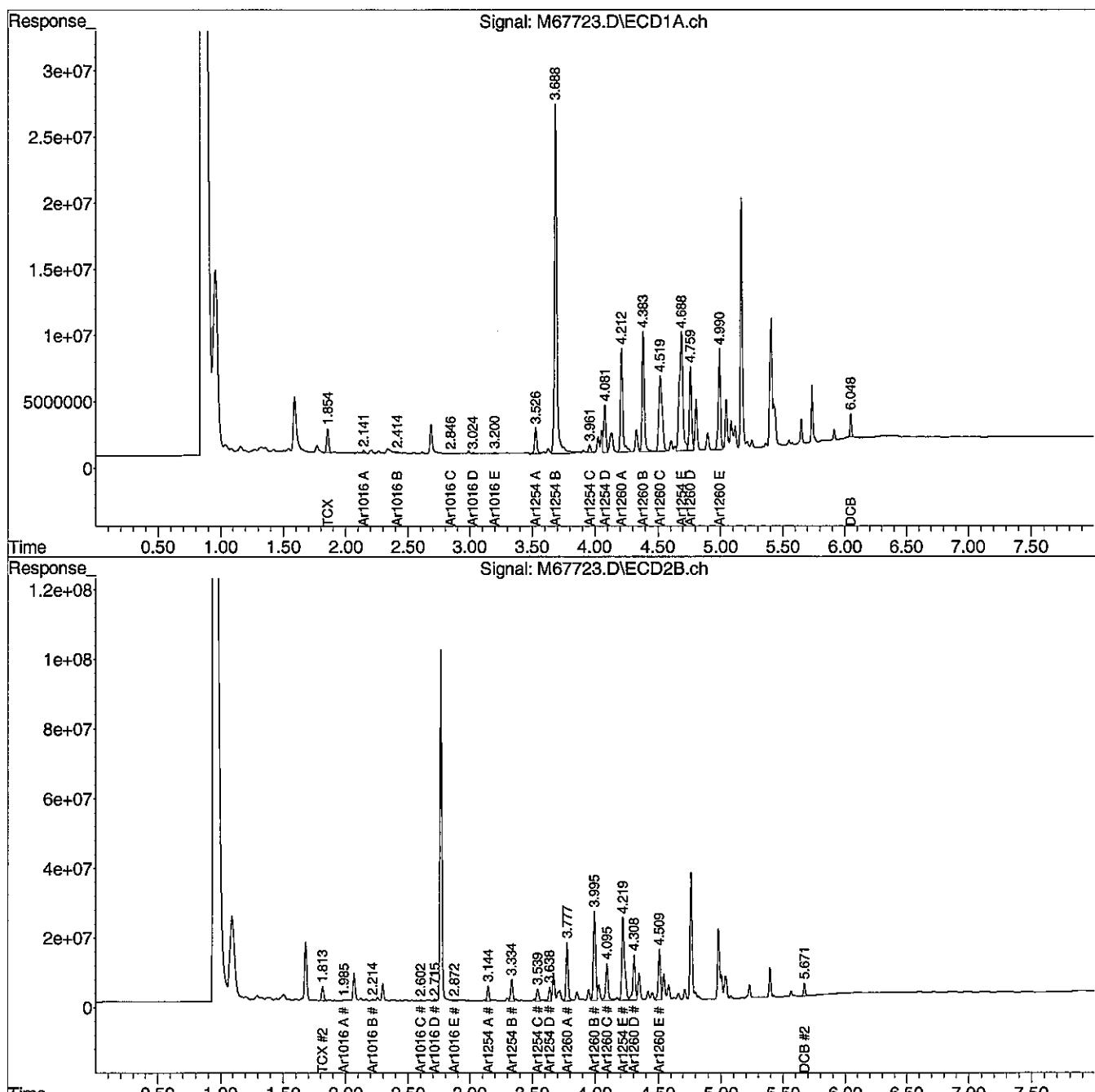
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67723.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:14 am
 Operator : JK
 Sample : 74887-8,1:10
 Misc : SOIL
 ALS Vial : 63 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:20 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-228 (8-10')

Lab Sample ID: 74887-9
Matrix: Solid
Percent Solid: 92
Dilution Factor: 52
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	1720	U
PCB-1221	1720	U
PCB-1232	1720	U
PCB-1242	1720	U
PCB-1248	1720	U
PCB-1254	1720	U
PCB-1260	1720	15500

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-9,1:50

Column ID: 0.25 mm

Data File: M67724.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 51.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	15248	15471	1.5	

Column to be used to flag RPD values greater than QC limit of 40%

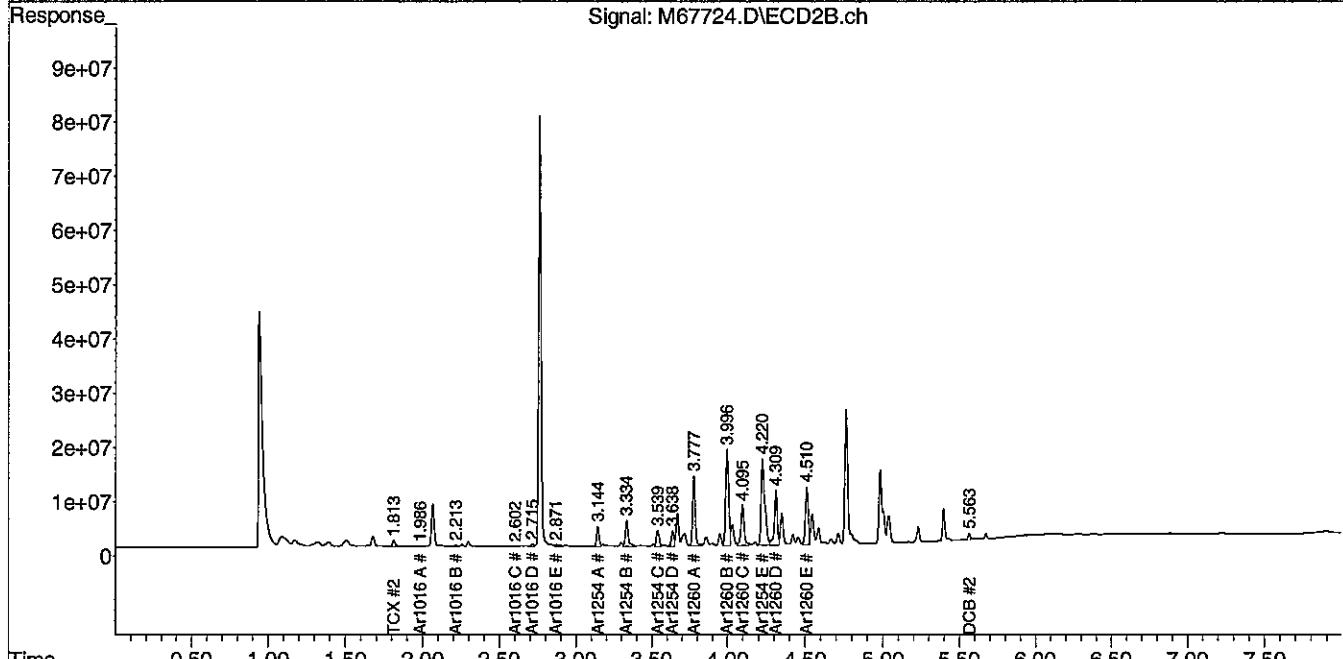
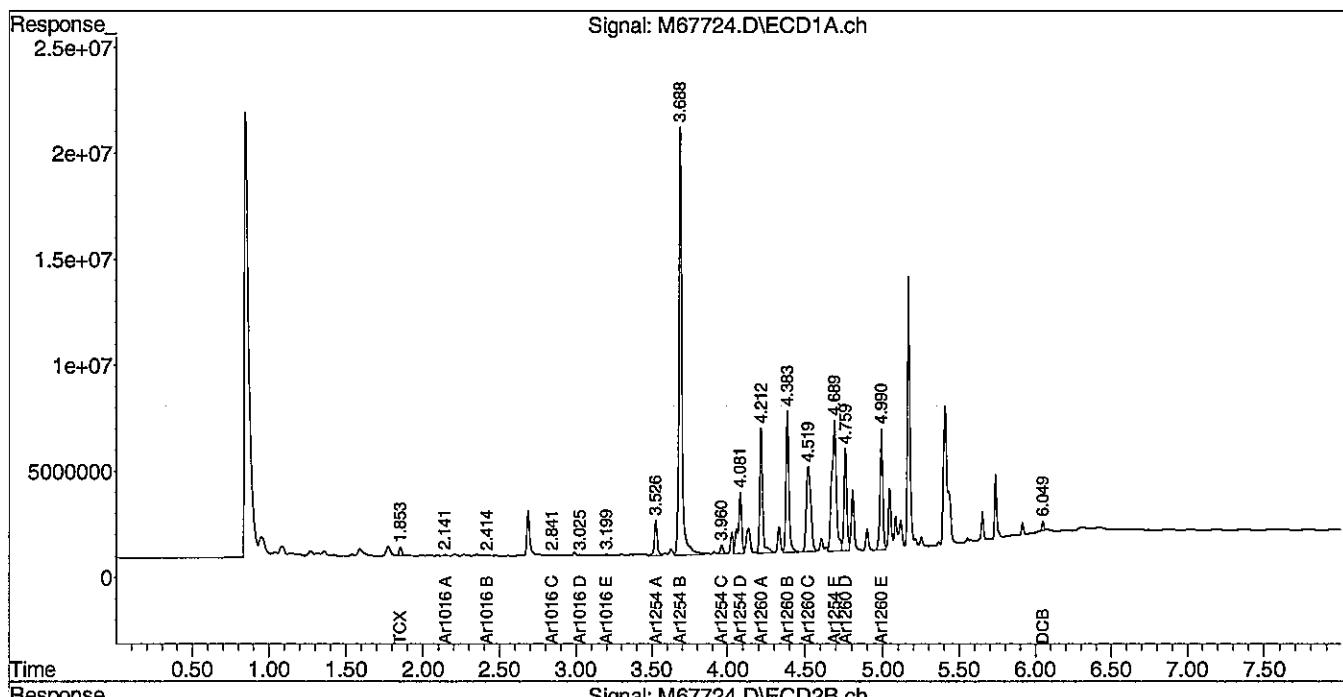
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67724.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:24 am
 Operator : JK
 Sample : 74887-9,1:50
 Misc : SOIL
 ALS Vial : 64 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:22 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-222 (0-2')

Lab Sample ID: 74887-10
Matrix: Solid
Percent Solid: 94
Dilution Factor: 10
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	330	U
PCB-1221	330	U
PCB-1232	330	U
PCB-1242	330	U
PCB-1248	330	U
PCB-1254	330	U
PCB-1260	330	2740

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	73	%
Decachlorobiphenyl	80	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-10,1:10

Column ID: 0.25 mm

Data File: M67725.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2744	2603	5.3	

Column to be used to flag RPD values greater than QC limit of 40%

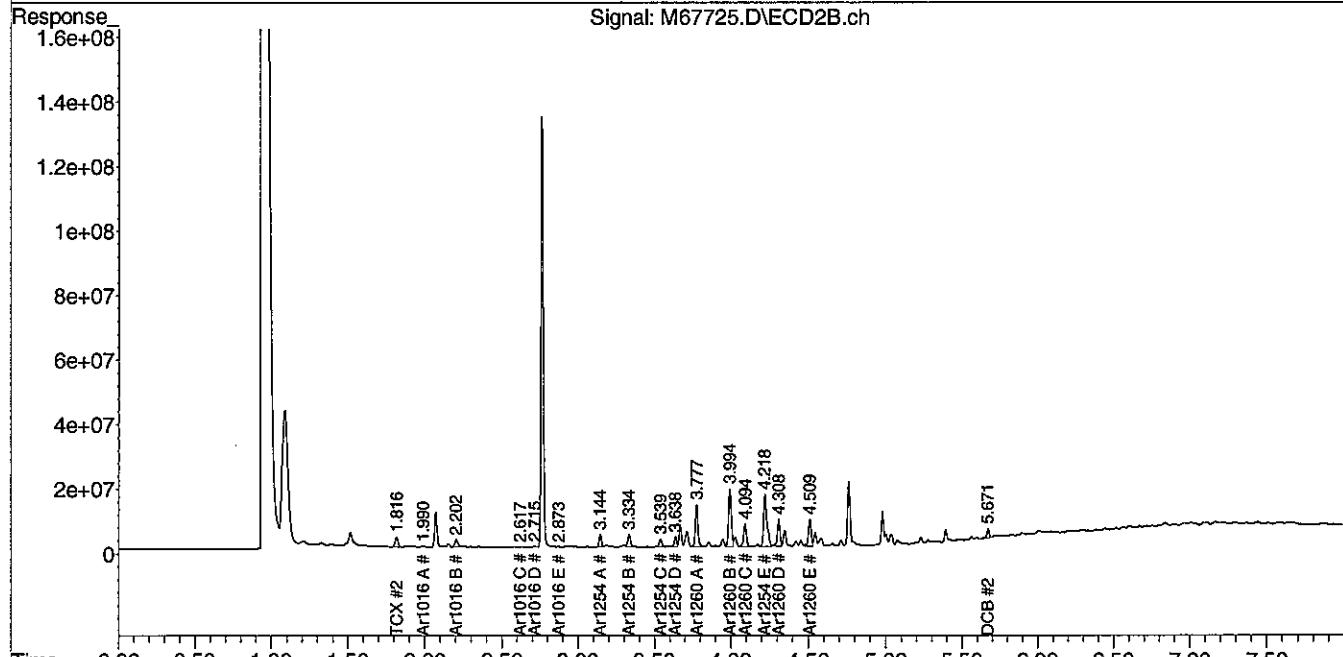
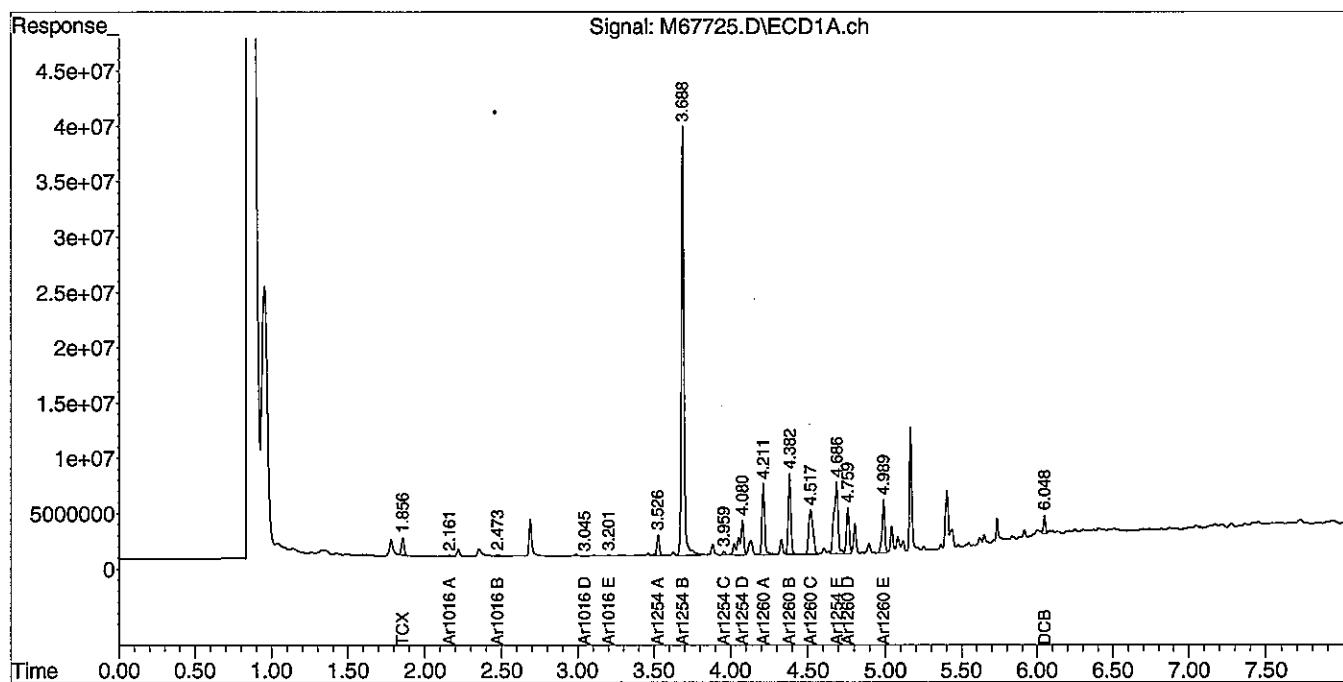
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67725.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:34 am
 Operator : JK
 Sample : 74887-10,1:10
 Misc : SOIL
 ALS Vial : 65 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:43:45 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-222 (2-4")

Lab Sample ID:	74887-11
Matrix:	Solid
Percent Solid:	92
Dilution Factor:	2.2
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	73	U
PCB-1221	73	U
PCB-1232	73	U
PCB-1242	73	U
PCB-1248	73	U
PCB-1254	73	U
PCB-1260	73	447

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	50	%
Decachlorobiphenyl	35*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * Surrogate recovery outside control limits. Secondary surrogate is in control. Sample was reanalyzed with similar results.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-11,1:2

Column ID: 0.25 mm

Data File: M67726.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	447	408	9.1	

Column to be used to flag RPD values greater than QC limit of 40%

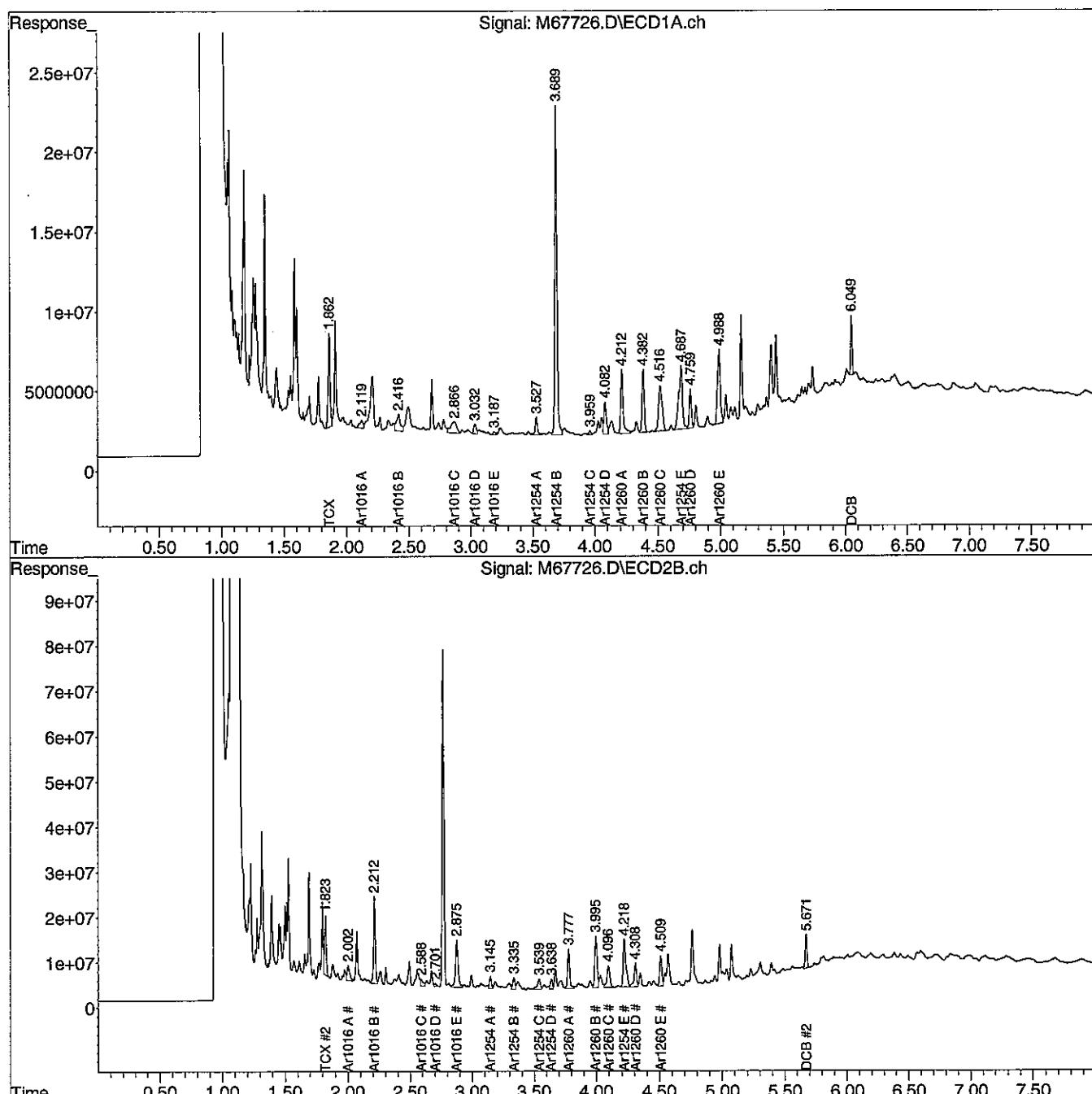
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67726.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:44 am
 Operator : JK
 Sample : 74887-11,1:2
 Misc : SOIL
 ALS Vial : 66 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:26 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-222 (4-6')

Lab Sample ID: 74887-12
Matrix: Solid
Percent Solid: 92
Dilution Factor: 2.1
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	69	U
PCB-1221	69	U
PCB-1232	69	U
PCB-1242	69	U
PCB-1248	69	U
PCB-1254	69	U
PCB-1260	69	823

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	86	%
Decachlorobiphenyl	77	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A.
 Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-12,1:5

Column ID: 0.25 mm

Data File: M67727.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	807	823	2.0	

Column to be used to flag RPD values greater than QC limit of 40%

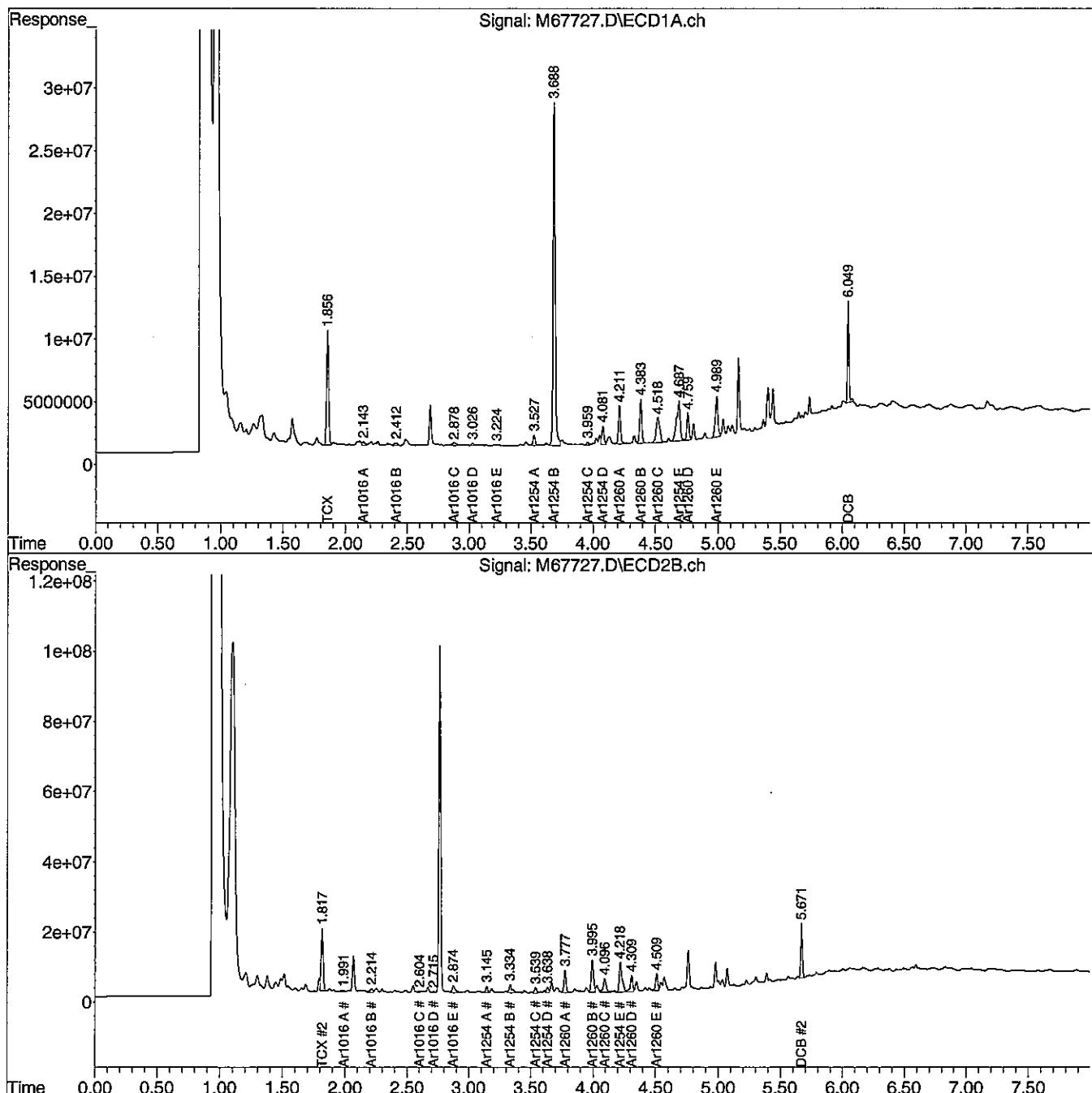
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67727.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 4:54 am
 Operator : JK
 Sample : 74887-12,1:2 (Sig #1); 74887-12,1:5 (Sig #2)
 Misc : SOIL
 ALS Vial : 67 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:28 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:41 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-222 (6-8')

Lab Sample ID: 74887-13
Matrix: Solid
Percent Solid: 95
Dilution Factor: 1.0
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	146

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	111	%
Decachlorobiphenyl	88	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-13

Column ID: 0.25 mm

Data File: M67733.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	116	146	23.0	

Column to be used to flag RPD values greater than QC limit of 40%

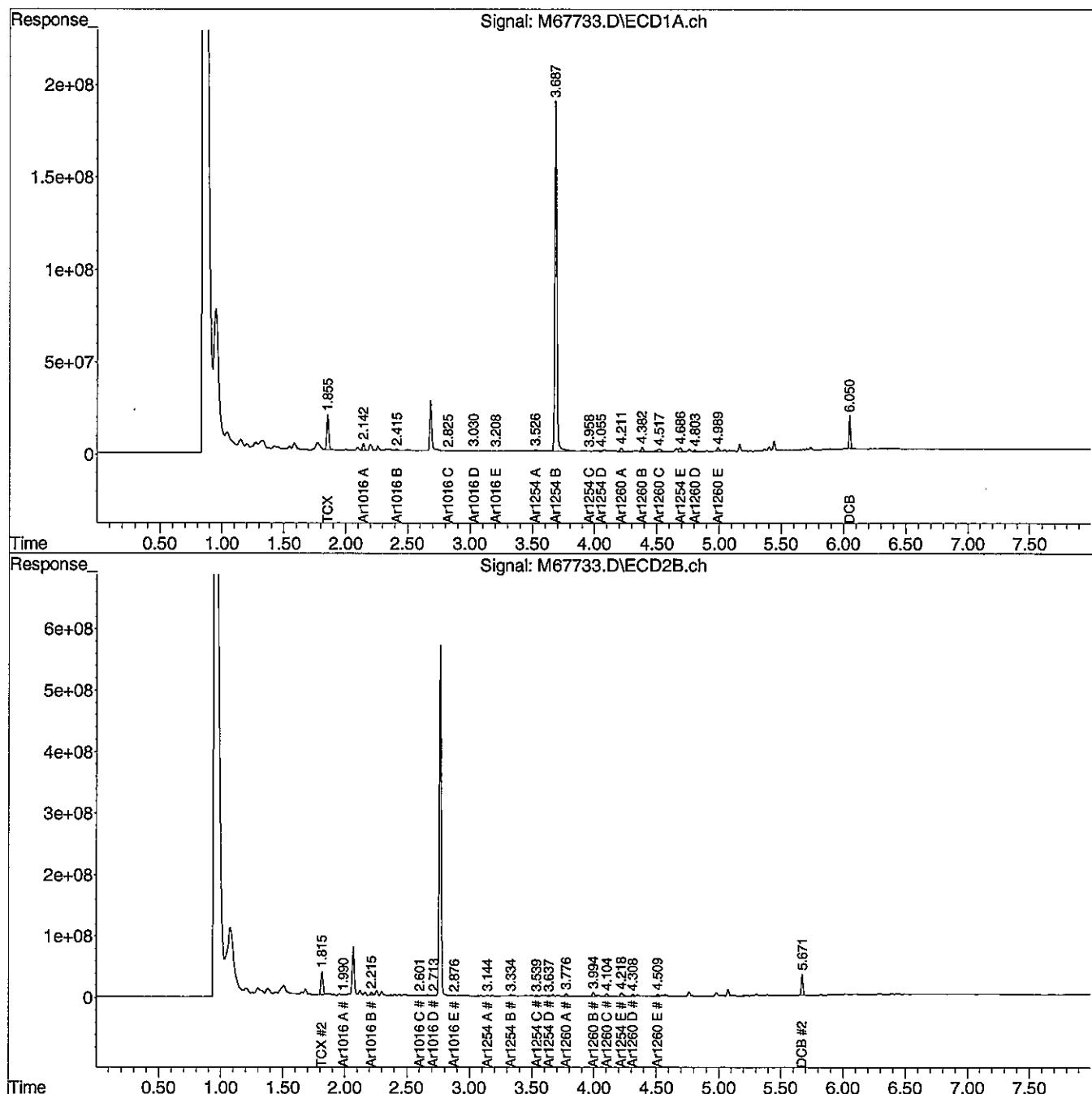
* Values outside QC limits

Comments: _____

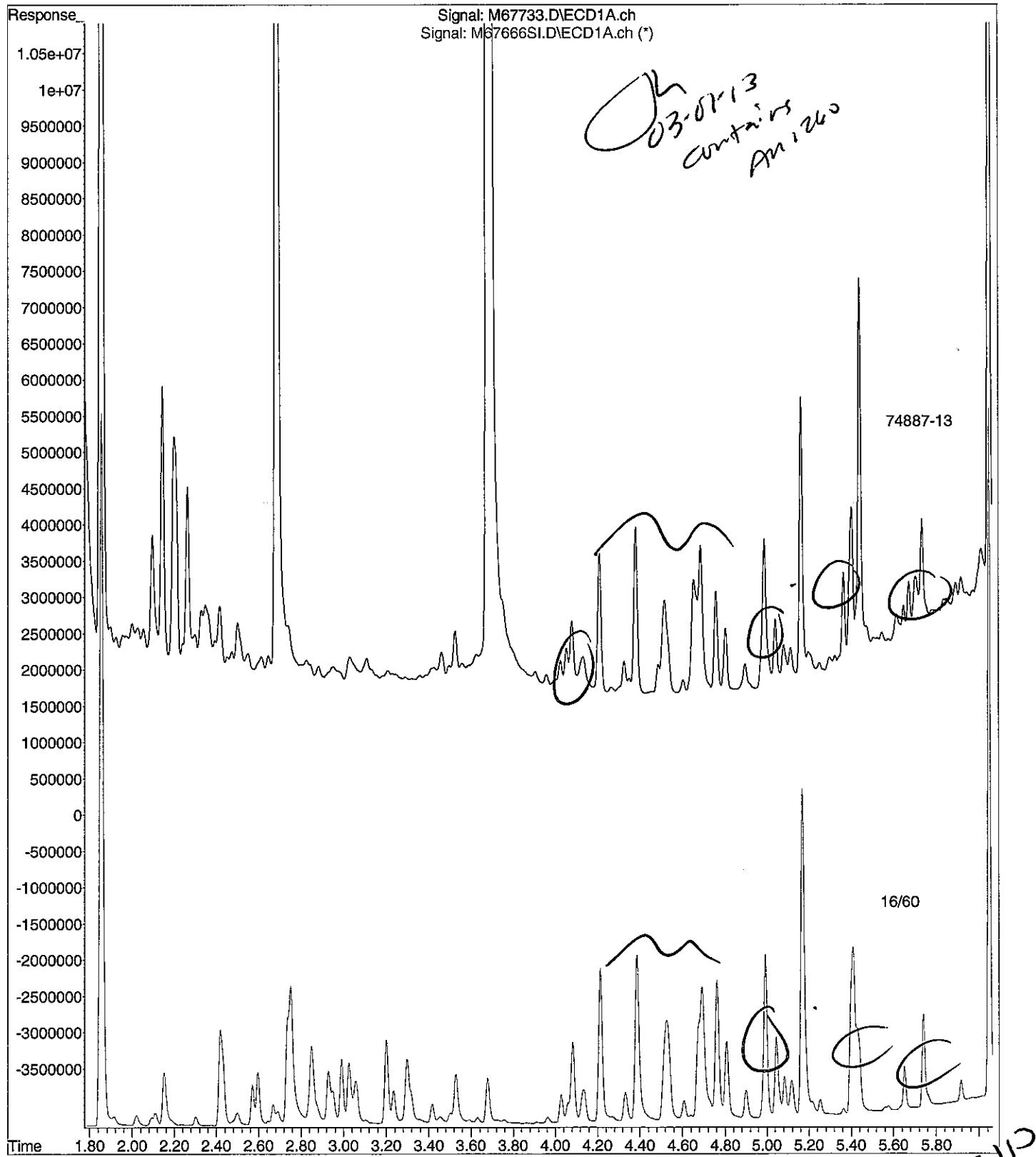
Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67733.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 5:54 am
 Operator : JK
 Sample : 74887-13
 Misc : SOIL
 ALS Vial : 68 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:27 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\022813-M\M67733.D
Operator : JK
Acquired : 1 Mar 2013 5:54 am using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74887-13
Misc Info : SOIL
Vial Number: 68



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-222 (8-10")

Lab Sample ID: 74887-14
Matrix: Solid
Percent Solid: 95
Dilution Factor: 1.0
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	237

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	83	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB Report

Authorized signature

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-14

Column ID: 0.25 mm

Data File: M67734.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	222	237	6.6	

Column to be used to flag RPD values greater than QC limit of 40%

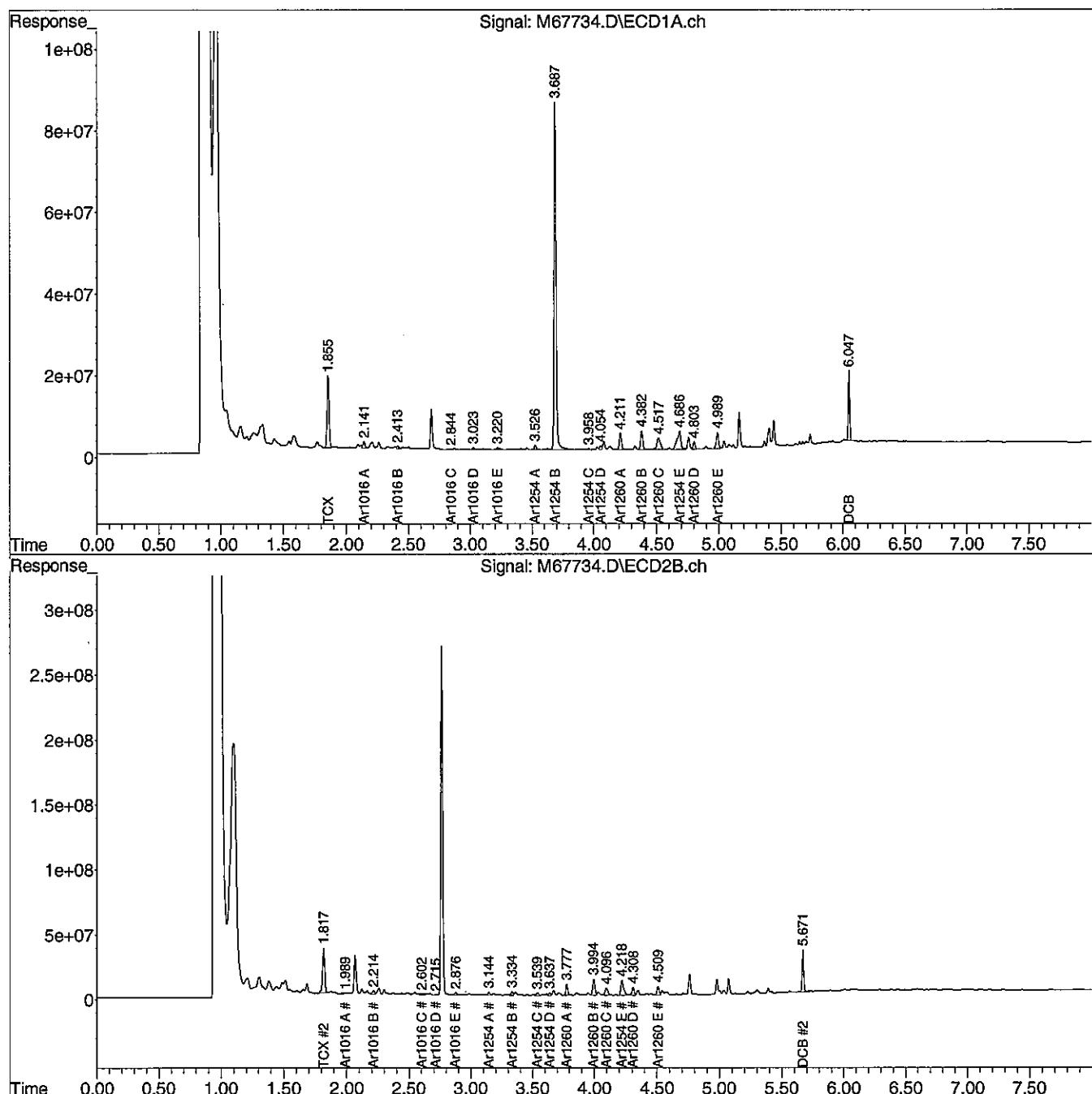
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67734.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 6:04 am
 Operator : JK
 Sample : 74887-14
 Misc : SOIL
 ALS Vial : 69 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:29 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-223 (0-2')

Lab Sample ID:	74887-15
Matrix:	Solid
Percent Solid:	95
Dilution Factor:	2.0
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/25/13
Analysis Date:	03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	66	U
PCB-1221	66	U
PCB-1232	66	U
PCB-1242	66	U
PCB-1248	66	U
PCB-1254	66	U
PCB-1260	66	133

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	98	%
Decachlorobiphenyl	79	%

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-15,1:2

Column ID: 0.25 mm

Data File: M67735.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	133	116	13.8	

Column to be used to flag RPD values greater than QC limit of 40%

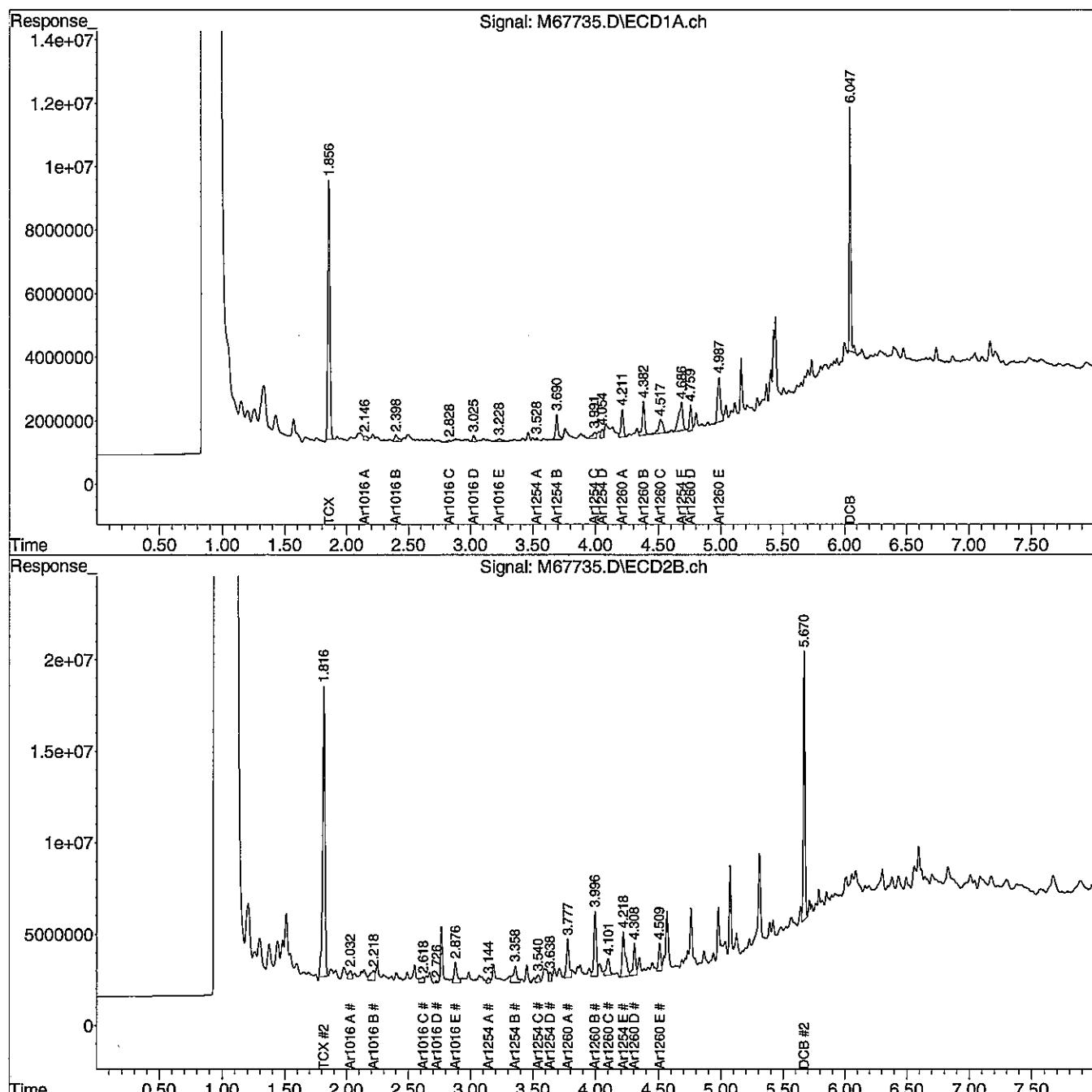
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67735.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 6:14 am
 Operator : JK
 Sample : 74887-15,1:2
 Misc : SOIL
 ALS Vial : 70 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:16:03 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-223 (2-4')

Lab Sample ID: 74887-16
Matrix: Solid
Percent Solid: 95
Dilution Factor: 488
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	16100	U
PCB-1221	16100	U
PCB-1232	16100	U
PCB-1242	16100	U
PCB-1248	16100	U
PCB-1254	16100	U
PCB-1260	16100	179000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M SDG: 74887
GC Column #1: STX-CLPesticides I Sample: 74887-16,1:500
Column ID: 0.25 mm Data File: M67736.D
GC Column #2: STX-CLPesticides II Dilution Factor: 487.9
Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	178568	168546	5.8	

Column to be used to flag RPD values greater than QC limit of 40%

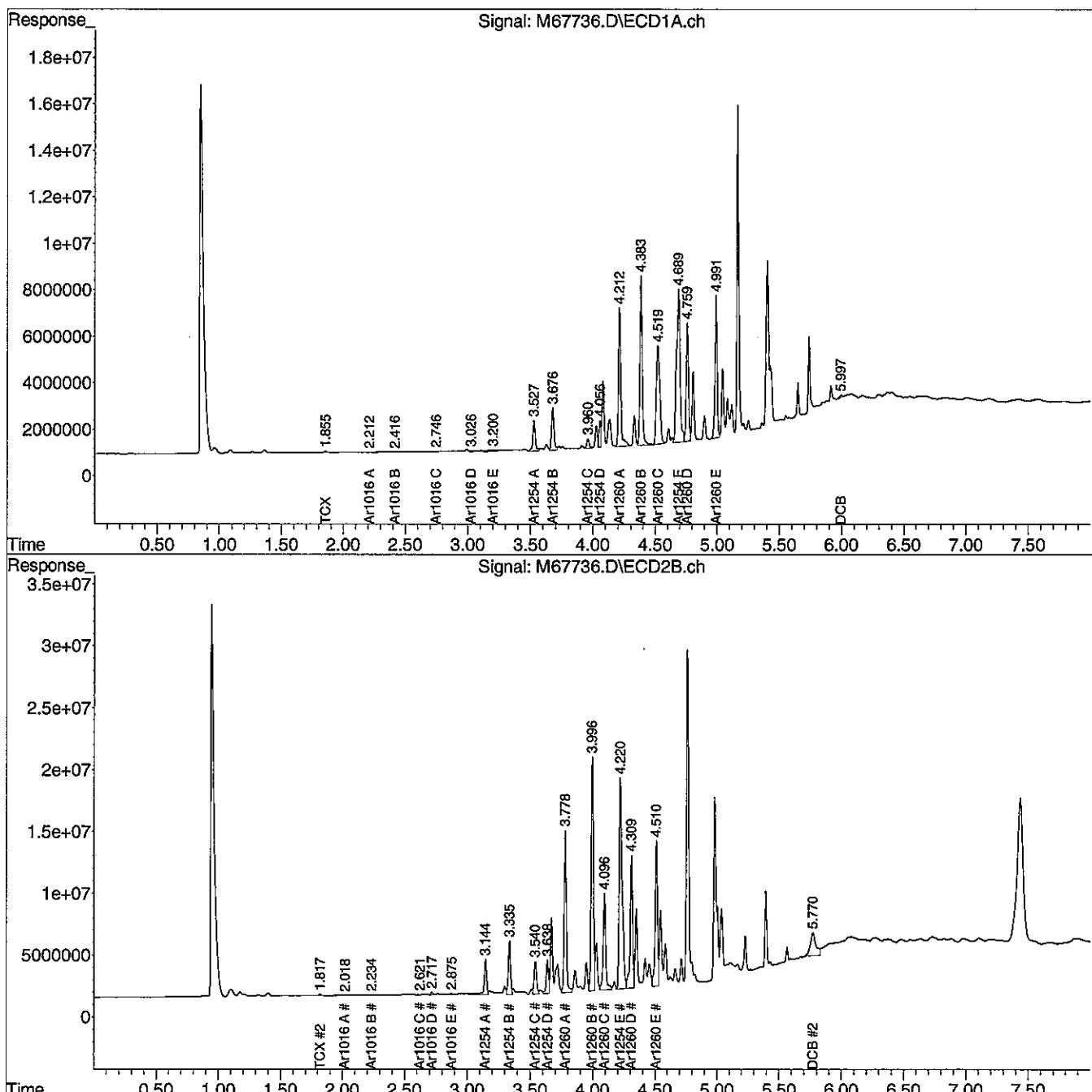
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67736.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 6:25 am
 Operator : JK
 Sample : 74887-16,1:500
 Misc : SOIL
 ALS Vial : 71 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:33 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-223 (4-6')

Lab Sample ID: 74887-17
Matrix: Solid
Percent Solid: 91
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	3920

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	110	%	
Decachlorobiphenyl	88	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-17,1:10

Column ID: 0.25 mm

Data File: M67737.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3924	3624	8.0	

Column to be used to flag RPD values greater than QC limit of 40%

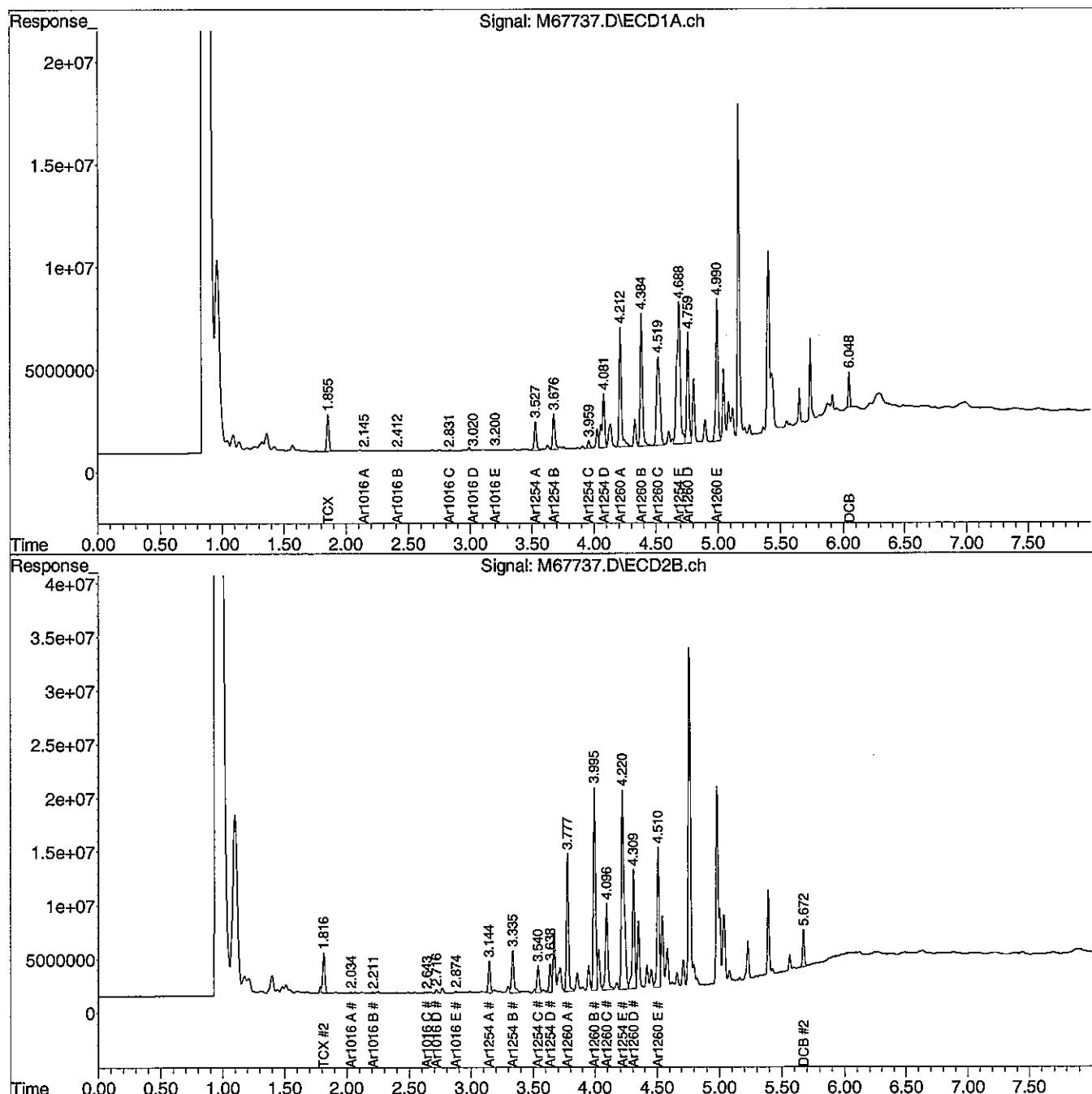
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67737.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 6:35 am
 Operator : JK
 Sample : 74887-17,1:10
 Misc : SOIL
 ALS Vial : 72 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:35 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: I2-3259.1
Field Sample ID: SE-SB-223 (6-8")

Lab Sample ID: 74887-18
Matrix: Solid
Percent Solid: 93
Dilution Factor: 11
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	363	U
PCB-1221	363	U
PCB-1232	363	U
PCB-1242	363	U
PCB-1248	363	U
PCB-1254	363	U
PCB-1260	363	3930

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	109	%
Decachlorobiphenyl	84	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-18,1:10

Column ID: 0.25 mm

Data File: M67738.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 10.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	3933	3450	13.1	

Column to be used to flag RPD values greater than QC limit of 40%

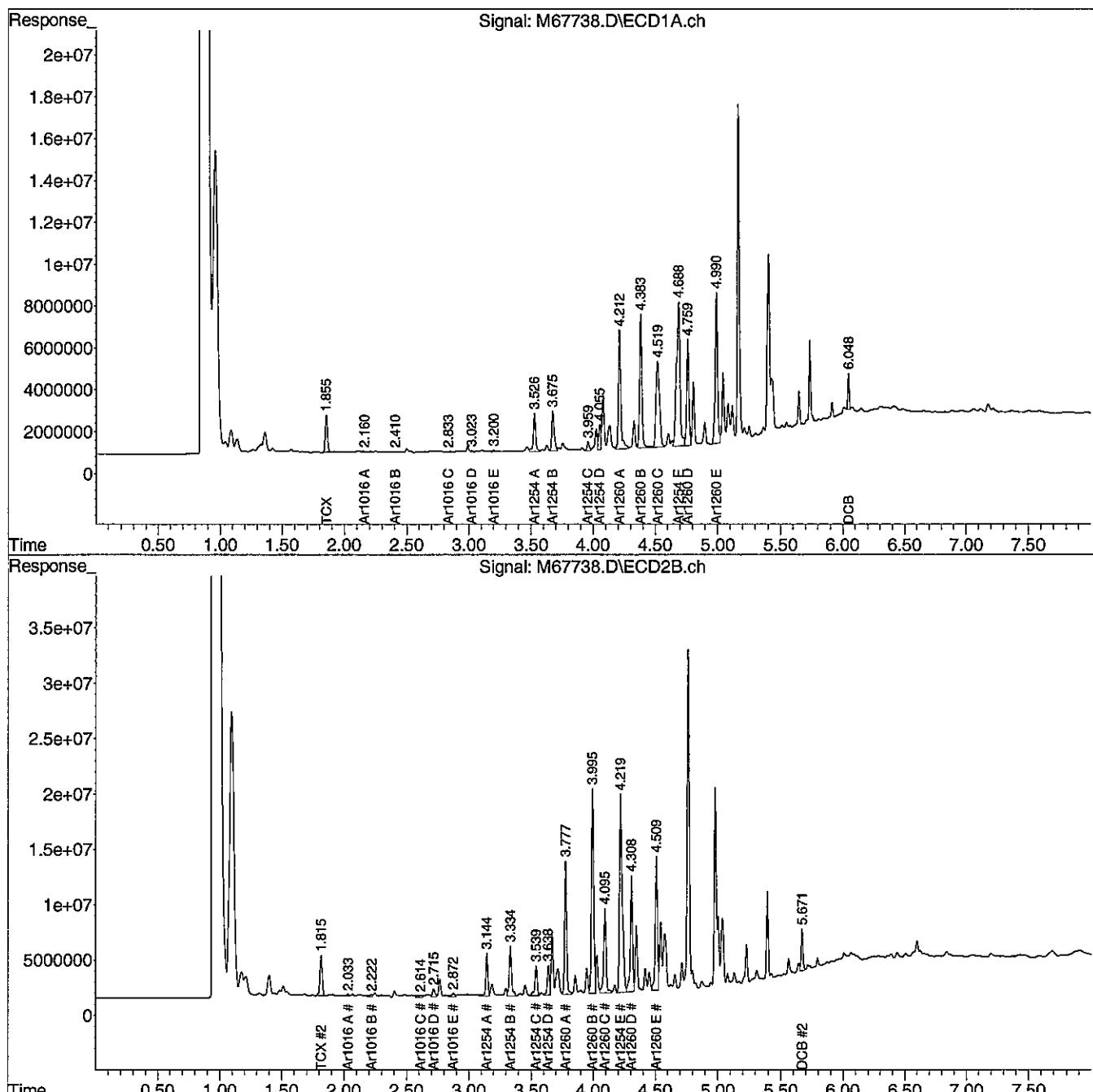
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67738.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 6:45 am
 Operator : JK
 Sample : 74887-18,1:10
 Misc : SOIL
 ALS Vial : 73 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:37 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-223 (8-10')

Lab Sample ID: 74887-19
Matrix: Solid
Percent Solid: 88
Dilution Factor: 22
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	726	U
PCB-1221	726	U
PCB-1232	726	U
PCB-1242	726	U
PCB-1248	726	U
PCB-1254	726	U
PCB-1260	726	8000

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-19,1:20

Column ID: 0.25 mm

Data File: M67741.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 22.3

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	7997	7183	10.7	

Column to be used to flag RPD values greater than QC limit of 40%

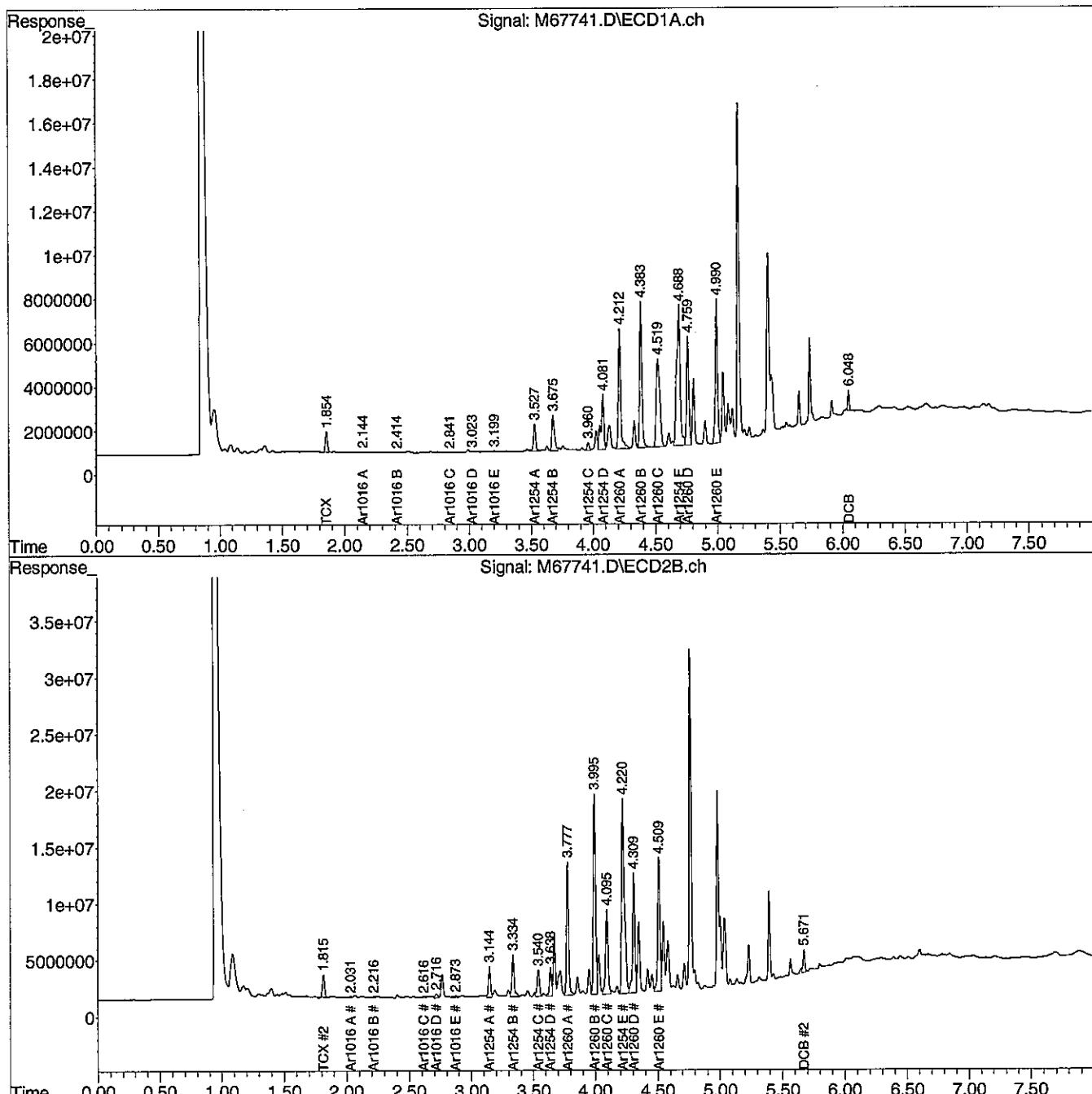
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67741.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 7:15 am
 Operator : JK
 Sample : 74887-19,1:20
 Misc : SOIL
 ALS Vial : 76 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:43 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-218 (0-2')

Lab Sample ID: 74887-20
Matrix: Solid
Percent Solid: 79
Dilution Factor: 2.4
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	79	U
PCB-1221	79	U
PCB-1232	79	U
PCB-1242	79	U
PCB-1248	79	U
PCB-1254	79	U
PCB-1260	79	328

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	100	%
Decachlorobiphenyl	100	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-20,1:2

Column ID: 0.25 mm

Data File: M67742.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.4

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	289	328	12.7	

Column to be used to flag RPD values greater than QC limit of 40%

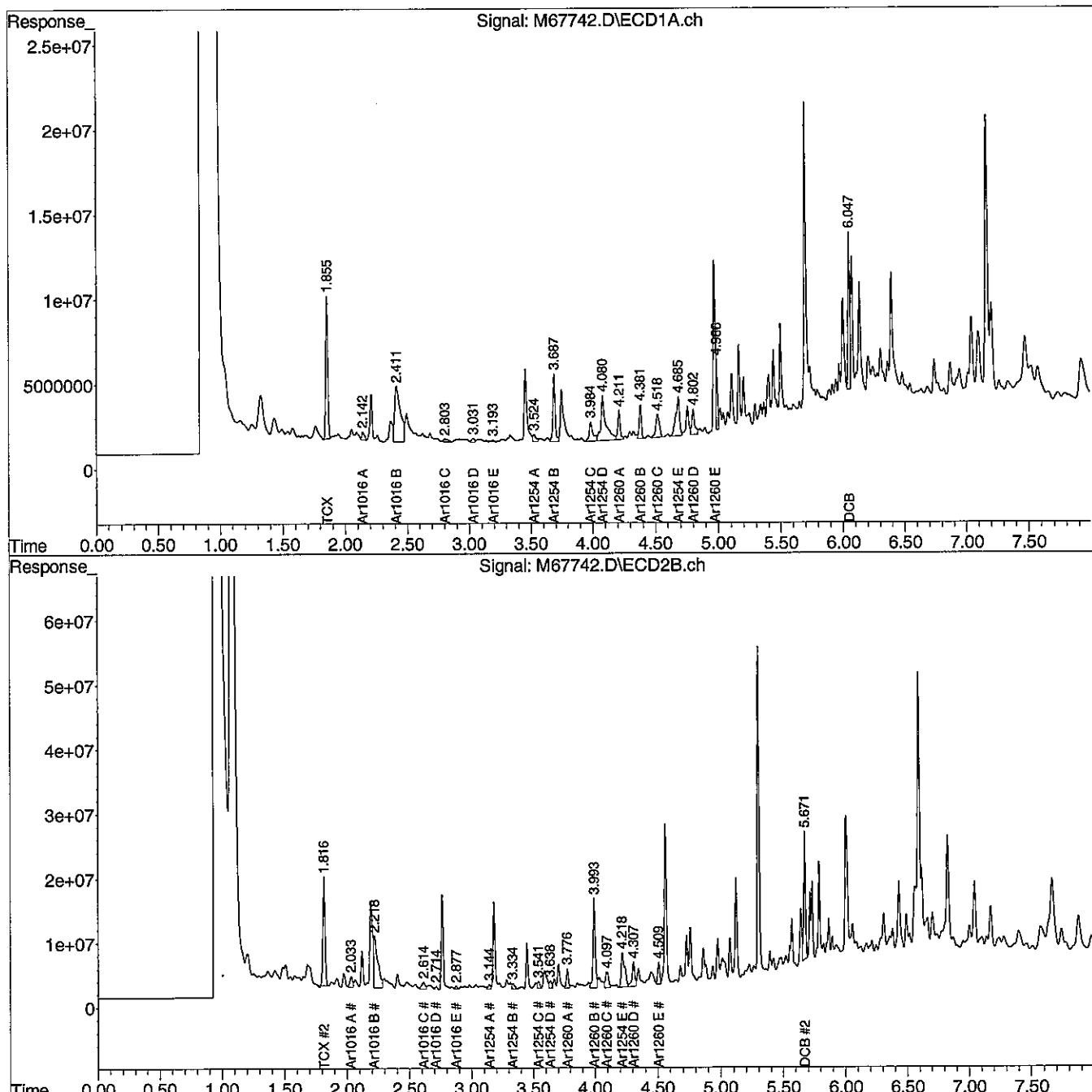
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67742.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 7:25 am
 Operator : JK
 Sample : 74887-20,1:2
 Misc : SOIL
 ALS Vial : 77 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:28:13 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: SE-SB-218 (2-4")

Lab Sample ID: 74887-21
Matrix: Solid
Percent Solid: 80
Dilution Factor: 1.2
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 02/26/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	40	U
PCB-1221	40	U
PCB-1232	40	U
PCB-1242	40	U
PCB-1248	40	U
PCB-1254	40	U
PCB-1260	40	696

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	88	%
Decachlorobiphenyl	83	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-21

Column ID: 0.25 mm

Data File: M67634.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD #
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)	
PCB 1260	696	546	24.1

Column to be used to flag RPD values greater than QC limit of 40%

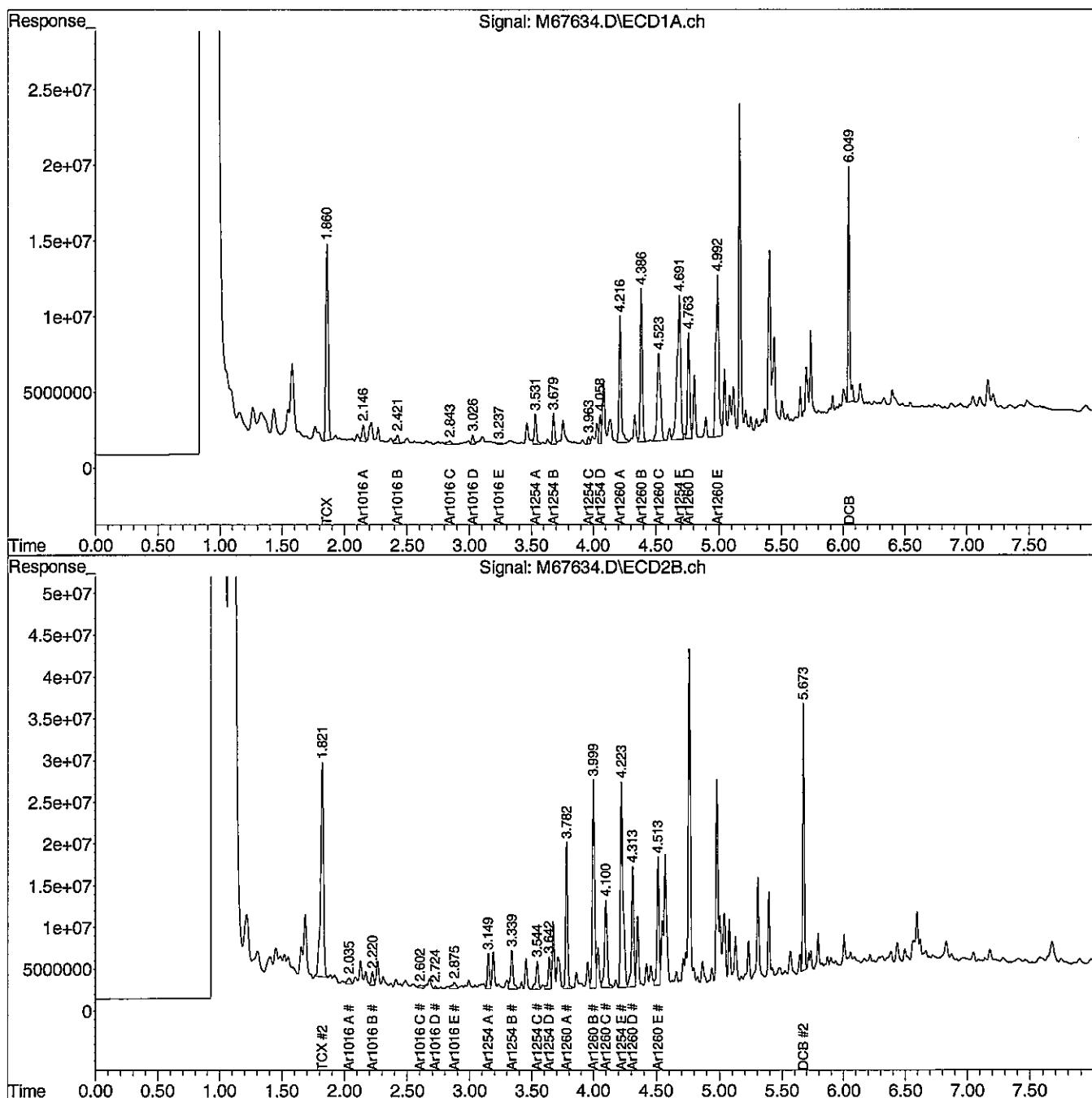
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67634.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 12:25 pm
 Operator : JK
 Sample : 74887-21
 Misc : SOIL
 ALS Vial : 12 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:52:05 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name:	MAINE ENERGY
Project Number:	12-3259.1
Field Sample ID:	SE-SB-231 (8-10')

Lab Sample ID:	74887-22
Matrix:	Solid
Percent Solid:	91
Dilution Factor:	5150
Collection Date:	02/20/13
Lab Receipt Date:	02/22/13
Extraction Date:	02/26/13
Analysis Date:	02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	170000	U
PCB-1221	170000	U
PCB-1232	170000	U
PCB-1242	170000	U
PCB-1248	170000	U
PCB-1254	170000	U
PCB-1260	170000	2330000

Surrogate Standard Recovery		
2,4,5,6-Tetrachloro-m-xylene	*	%
Decachlorobiphenyl	*	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.
 * The surrogates were diluted out.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74887

GC Column #1: STX-CLPesticides I

Sample: 74887-22,1:5000

Column ID: 0.25 mm

Data File: M67639.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5144.6

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2332584	2204134	5.7	

Column to be used to flag RPD values greater than QC limit of 40%

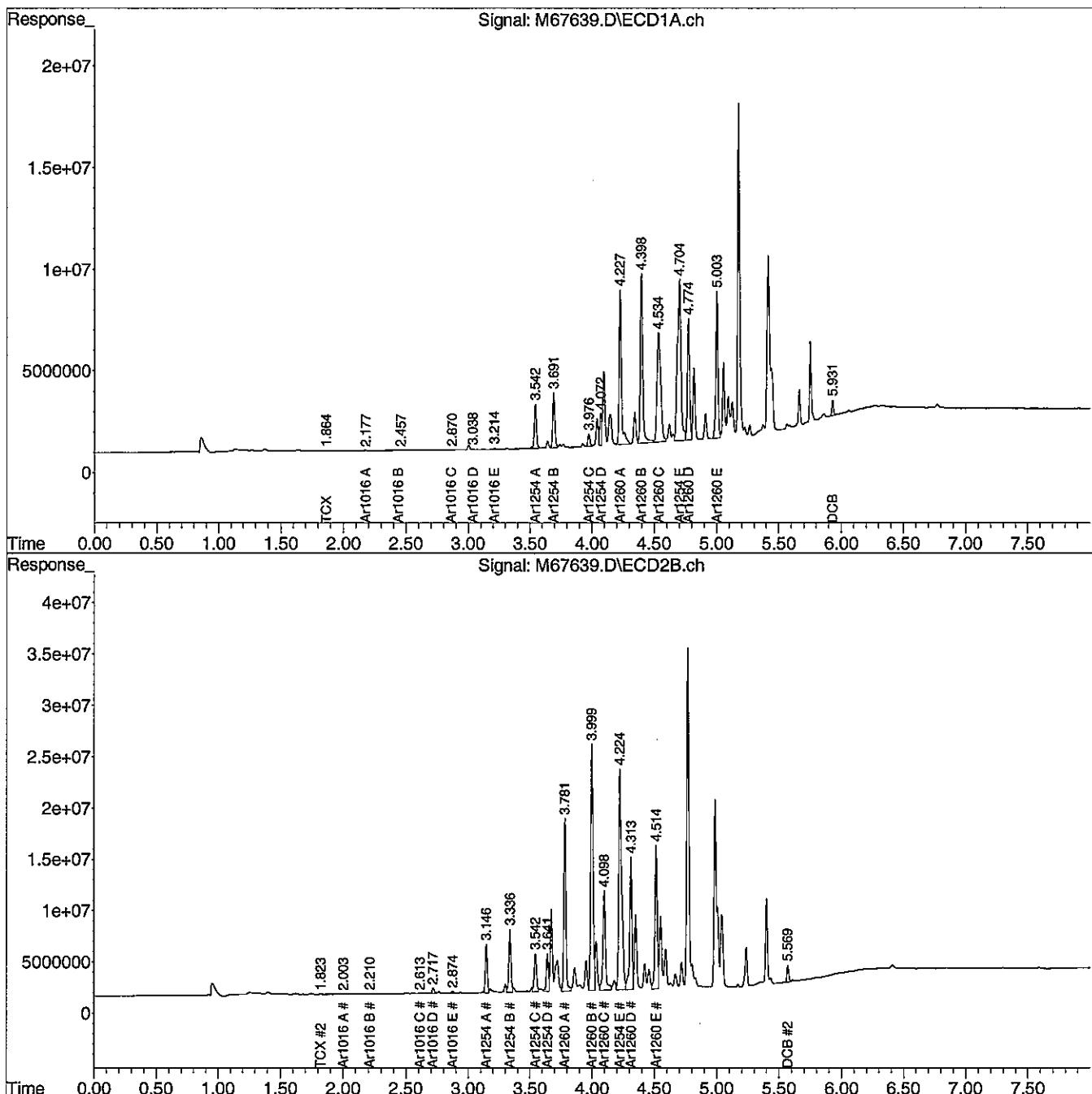
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67639.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 28 Feb 2013 1:34 pm
 Operator : JK
 Sample : 74887-22,1:5000
 Misc : SOIL
 ALS Vial : 17 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Feb 28 13:52:15 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:13 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



PCB
QC FORMS

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February 28, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022613PSOX
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/26/13
Analysis Date: 02/28/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	77	%
Decachlorobiphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

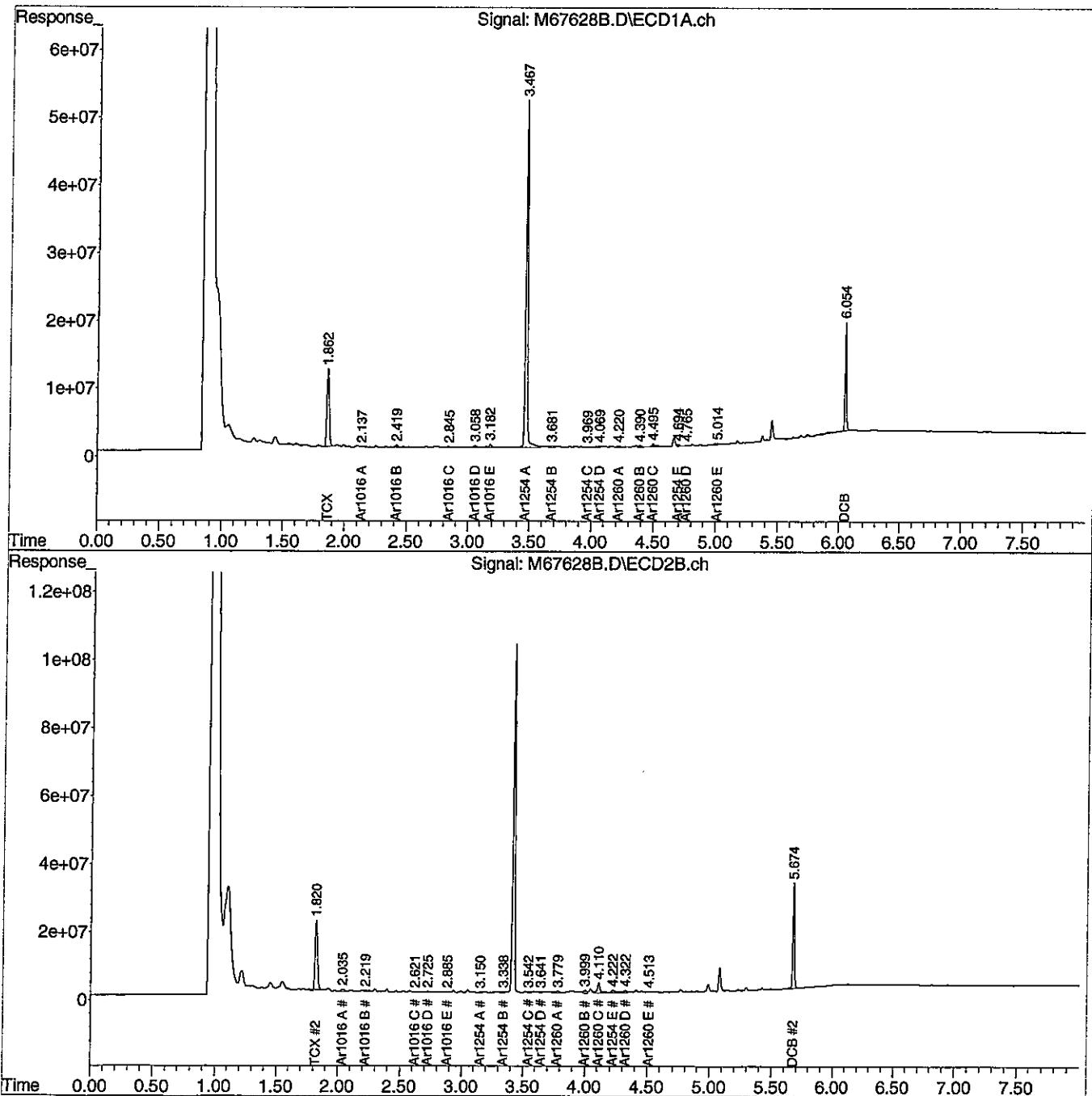
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022813-M\
Data File : M67628B.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 28 Feb 2013 11:24 am
Operator : JK
Sample : B022613PSOX
Misc : SOIL
ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
Integration File signal 2: events2.e
Quant Time: Feb 28 13:51:53 2013
Quant Method : C:\msdchem\1\METHODS\PCB020513.M
Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
QLast Update : Thu Feb 28 11:22:11 2013
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 2 uL
Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



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March 1, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022513PSOX2
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	65	%
Decachlorobiphenyl	81	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

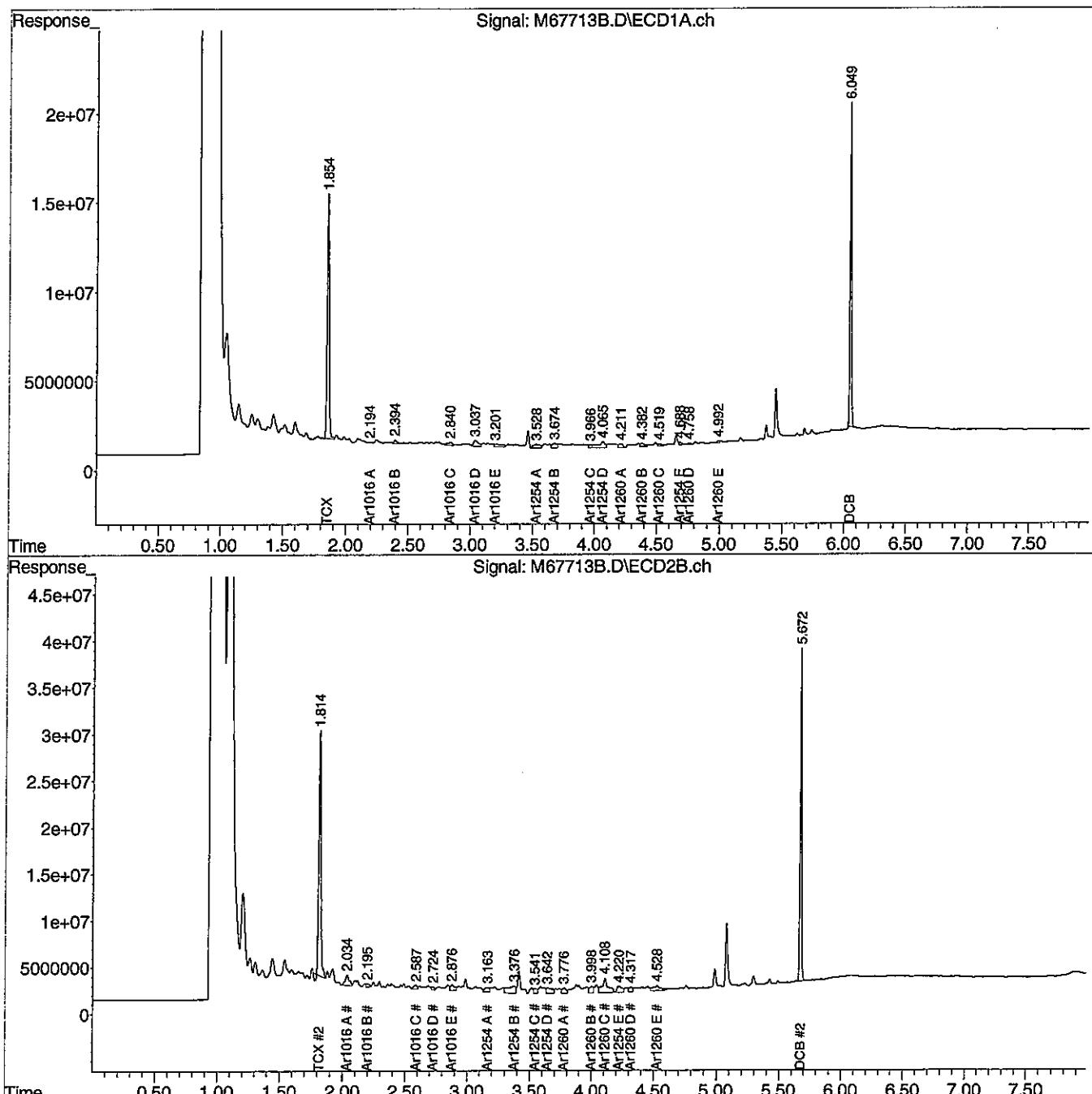
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67713B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 2:33 am
 Operator : JK
 Sample : B022513PSOX2
 Misc : SOIL
 ALS Vial : 53 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 10:33:00 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Fri Mar 01 08:56:39 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



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March 1, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: MAINE ENERGY
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B022513PSOX2 RR
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 02/25/13
Analysis Date: 03/01/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	81	%
Decachlorobiphenyl	84	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

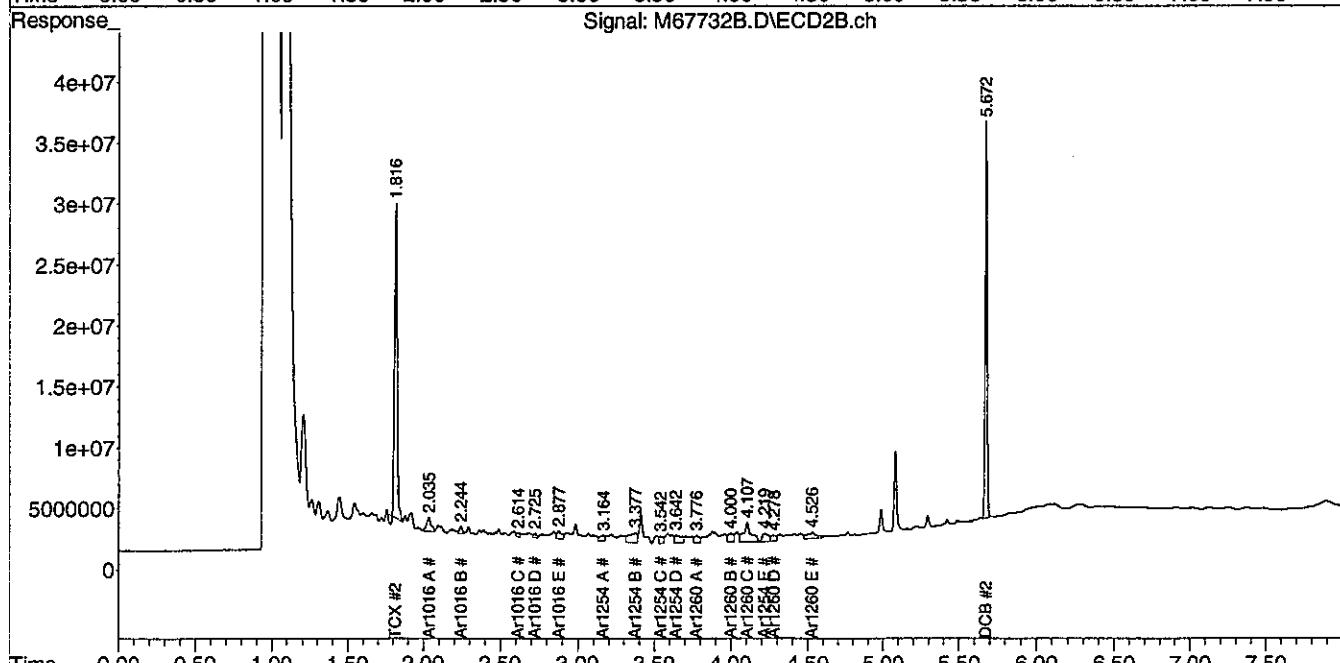
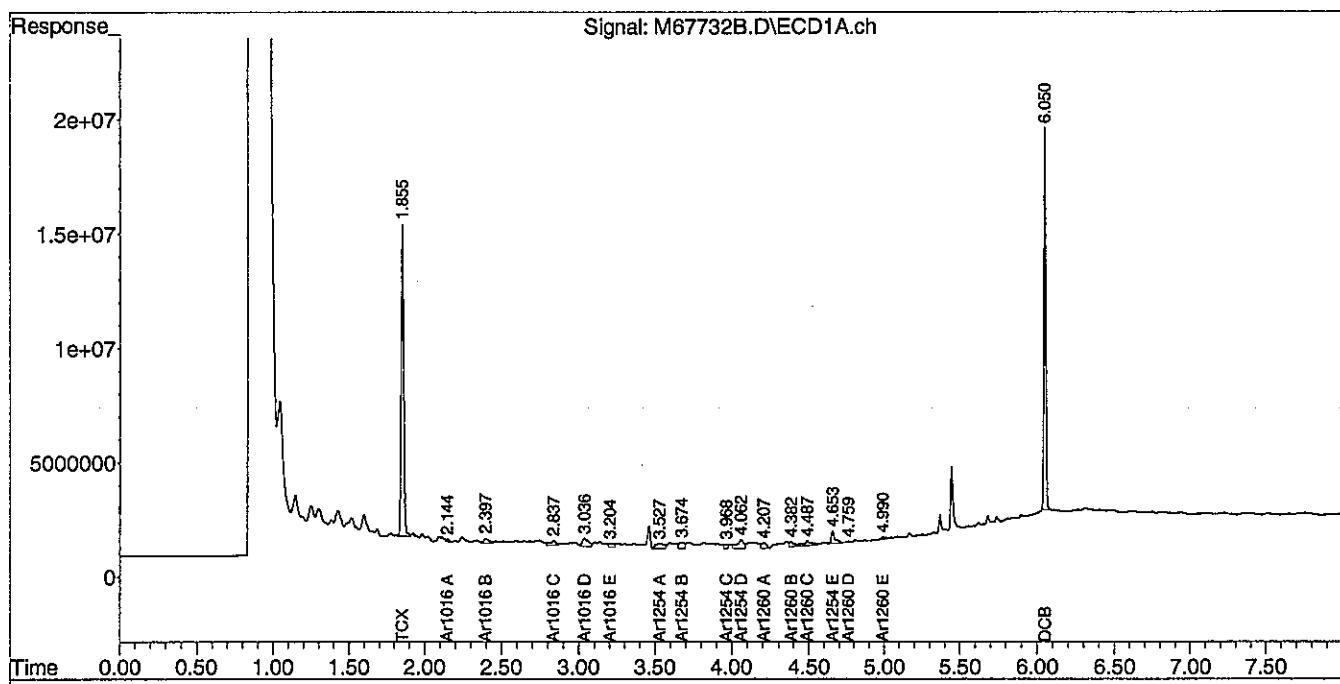
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\022813-M\
 Data File : M67732B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 1 Mar 2013 5:44 am
 Operator : JK
 Sample : B022513PSOX2,RR
 Misc : SOIL
 ALS Vial : 53 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 01 11:12:25 2013
 Quant Method : C:\msdchem\1\METHODS\PCB020513.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Thu Feb 28 11:22:11 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74887

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside QC limits

D System Monitoring Compound diluted out

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SYSTEM MONITORING COMPOUNDS
SUMMARY**

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PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022613PSOX

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: L022613PSOX

Spike duplicate: LD022613PSOX

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP	RPD	#	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#		
PCB 1016	200	200	65	140	30	0	181	90		195	97		7.5	
PCB 1260	200	200	60	130	30	0	182	91		187	94		2.9	
PCB 1016 #2	200	200	65	140	30	0	167	83		179	90		7.3	
PCB 1260 #2	200	200	60	130	30	0	192	96		189	94		1.7	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

SDG:

Non-spiked sample: B022513PSOX2

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

Spike: L022513PSOX2

Spike duplicate: LD022513PSOX2

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP			
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	#
PCB 1016	200	200	65	140	30	0	186	93		172	86		7.9	
PCB 1260	200	200	60	130	30	0	173	87		178	89		2.6	
PCB 1016 #2	200	200	65	140	30	0	183	91		176	88		3.8	
PCB 1260 #2	200	200	60	130	30	0	208	104		190	95		9.0	

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG: 74887

Non-spiked sample: 74887-18,1:10

Spike: 74887-18,MS,1:10

Spike duplicate: 74887-18,MSD,1:11

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE	SPIKE DUP	SPIKE DUP	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC #	RESULT (ug/kg)	% REC #	RPD #
PCB 1016	204	213	65	140	30	0	276	135	274	129	1.0
PCB 1260	204	213	60	130	30	3933	4767	408 *	3731	-95 *	24.4
PCB 1016 #2	204	213	65	140	30	0	221	108	232	109	5.2
PCB 1260 #2	204	213	60	130	30	3450	3013	-214 *	3389	-29 *	11.7

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form

 Project Name: MAIN ENERGY		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax		For Analytics Use Only	
Project#: 12-3359.1		Preservation Code: Preservation Key: A = HCl B = 4 C C = Unpres D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other		Circle and/or Write Required Analysis Followed by Preservation Code Please initial in preservation code here		Samples were: 1) Shipped <input checked="" type="checkbox"/> Hand-delivered <input type="checkbox"/> 2) Temperature (°C): 40 3) Received in good condition: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No 4) pH checked by: KCJ 02/22/13 5) Labels checked by:	
Company: SMC						Matrix Key: C = Concrete WP = Wipe SW = Wastewater SW = Surface Water E = Extract	
Report to: JOHN CRESSEY						GW = Groundwater DW = Drinking Water S = Soil / Sludge O = Oil X = Other	
Address: 64 Main St. Lewiston ME 04240							
Phone: (607) 745-5007							
Quote #:							
PO# (if required): 12-3359.1							
Sample Identification		Sample Date	Sample Time	Sample	Time	Matrix	No. of Containers checked
SE-13-227 (0-3)		2-20	1345	N		S	1
SE-13-227 (0-1)		2-20	1350	N		S	1
SE-13-227 (4-6)		2-20	1351	N		S	1
SE-13-227 (6-8)		2-20	1354	N		S	1
SE-13-228 (0-3)		2-20	1400	N		S	1
SE-13-228 (0-1)		2-20	1404	N		S	1
SE-13-228 (4-6)		2-20	1408	N		S	1
SE-13-228 (6-8)		2-20	1414	N		S	1
SE-13-228 (8-10)		2-20	1416	N		S	1
SE-13-228 (0-3)		2-20	1440	N		S	1
** List requested metals here +Fee may apply							
Comments, Additional Analyses, or Special Instructions: Email Results to: JOHN CRESSEY							
Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 4 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> Standard (8-10 business days)							
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.							
*Fee may apply; lab approval required							
Sampler Name (Print): JOHN CRESSEY		Date: 2-21 Date: 02/03		Time: 9:00 Time: 8:50		Received By: General Counsel Received By: John Cresssey	
Relinquished By Sampler:							
Relinquished By:							
Relinquished By:							

Chain Of Custody Form

 Maine Enviro LLC environmental laboratory		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax	
Project Name: MAINE ENVIRONMENTAL Project #: 12-3259-1					
Company: Symonit Report to: John Cressy Address: 47 Main St. City: Concord, ME 04401 Phone: (207) 765-4507 Quote #: PO# (if required): 12-3259-1					
Preservation Code: Please圈出 in preservation code here					
Preservation Key: A = HCl B = 4°C C = Unpres. D = MeOH E = HNO3 F = H2SO4 G = Hexane H = Other					
Sample Identification					
Sample Date	Sample Time	Field Filtered Y or N	VOC: 8260 524.2 624	SVOC: 8270 625 PAH only SIM	Pesticides: 8081 608
SE-38-222 (2-4)	2-20	X	X	X	PCB: 8082 508 Soxhlet Y or N
SE-38-222 (4-6)	2-20	X	X	X	TPH: 8015 (Gas Range)
SE-38-222 (6-8)	2-20	X	X	X	TPH: 8015 (Diesel Range)
SE-38-222 (8-10)	2-20	X	X	X	TPH: Full or Ranges only
SE-38-223 (0-2)	2-20	X	X	X	VPH: Full or Ranges only TEPH
SE-38-223 (2-4)	2-20	X	X	X	Metals: Full or Ranges only
SE-38-223 (4-6)	2-20	X	X	X	GW = Groundwater
SE-38-223 (6-8)	2-20	X	X	X	DW = Drinking Water
SE-38-223 (8-10)	2-20	X	X	X	WW = Wastewater
SE-38-223 (0-2)	2-20	X	X	X	SW = Surface Water
SE-38-223 (2-4)	2-20	X	X	X	E = Extract
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
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SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
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SE-38-223 (2-4)	2-20	X	X	X	
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SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
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SE-38-223 (8-10)	2-20	X	X	X	
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SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
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SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
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SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)	2-20	X	X	X	
SE-38-223 (2-4)	2-20	X	X	X	
SE-38-223 (4-6)	2-20	X	X	X	
SE-38-223 (6-8)	2-20	X	X	X	
SE-38-223 (8-10)	2-20	X	X	X	
SE-38-223 (0-2)</td					

Chain Of Custody Form

Analytics environmental Laboratory LLC		195 Commerce Way, Suite E Portsmouth, NH 03801 (800) 929-9906		(603) 436-5111 (603) 430-2151 Fax																					
For Analytics Use Only																									
Samples were: 1) Shipped or hand-delivered <input checked="" type="checkbox"/> <input type="checkbox"/> 2) Temperature (°C): <input type="text" value="40"/> 3) Received in good condition: <input checked="" type="checkbox"/> Yes <input type="checkbox"/> No 4) pH checked by: <input type="text" value="423072243"/> 5) Labels checked by:																									
Matrix Key: C = Concrete WP = Wine SW = Wastewater SW = Surface Water E = Extract																									
Matrix: <input type="text" value="S"/> No. of Containers checked: <input type="text" value="1"/> pH: <input type="text" value="74887-21"/>																									
Circle and/or Write Required Analysis Followed by Preservation Code Please fill in preservation code here																									
* * * * *																									
Metals: RCRA PPs13 TAL23 Other * * * * *																									
VPH: Full or Ranges only TEPH																									
EPH: Full or Ranges only S100M ME125																									
TPB: 8015 (Diesel Range) ME217																									
TPB: 8015 (Gas Range) ME217																									
PCB: 8082 608 Solvent only N																									
VOC: 8260 524.2 624																									
SVOC: 8270 625 Path only S1A																									
Pesticides: 8081 608																									
SVOC: 8270 625 Path only S1A																									
** List requested metals here																									
Project Requirements: * Fee may apply																									
Report Type: <table border="1" style="margin-left: auto; margin-right: auto;"> <tr> <td><input type="checkbox"/> MCP*</td> <td><input type="checkbox"/> Level I*</td> <td><input type="checkbox"/> NH</td> <td><input type="checkbox"/> State:</td> </tr> <tr> <td><input type="checkbox"/> CTRCP*</td> <td><input type="checkbox"/> Level II*</td> <td><input type="checkbox"/> MA</td> <td><input type="checkbox"/> State Standard:</td> </tr> <tr> <td><input type="checkbox"/> DOD*</td> <td><input type="checkbox"/> Level IV*</td> <td><input checked="" type="checkbox"/> ME</td> <td><input type="checkbox"/> (eg. S-1 or GW-1)</td> </tr> <tr> <td><input checked="" type="checkbox"/> Standard</td> <td><input type="checkbox"/> RI</td> <td><input type="checkbox"/> CT</td> <td><input type="checkbox"/> EDD Required: Y* <input checked="" type="checkbox"/> N</td> </tr> <tr> <td colspan="2"></td> <td><input type="checkbox"/> Other</td> <td>Type: <input type="text" value=""/></td> </tr> </table>						<input type="checkbox"/> MCP*	<input type="checkbox"/> Level I*	<input type="checkbox"/> NH	<input type="checkbox"/> State:	<input type="checkbox"/> CTRCP*	<input type="checkbox"/> Level II*	<input type="checkbox"/> MA	<input type="checkbox"/> State Standard:	<input type="checkbox"/> DOD*	<input type="checkbox"/> Level IV*	<input checked="" type="checkbox"/> ME	<input type="checkbox"/> (eg. S-1 or GW-1)	<input checked="" type="checkbox"/> Standard	<input type="checkbox"/> RI	<input type="checkbox"/> CT	<input type="checkbox"/> EDD Required: Y* <input checked="" type="checkbox"/> N			<input type="checkbox"/> Other	Type: <input type="text" value=""/>
<input type="checkbox"/> MCP*	<input type="checkbox"/> Level I*	<input type="checkbox"/> NH	<input type="checkbox"/> State:																						
<input type="checkbox"/> CTRCP*	<input type="checkbox"/> Level II*	<input type="checkbox"/> MA	<input type="checkbox"/> State Standard:																						
<input type="checkbox"/> DOD*	<input type="checkbox"/> Level IV*	<input checked="" type="checkbox"/> ME	<input type="checkbox"/> (eg. S-1 or GW-1)																						
<input checked="" type="checkbox"/> Standard	<input type="checkbox"/> RI	<input type="checkbox"/> CT	<input type="checkbox"/> EDD Required: Y* <input checked="" type="checkbox"/> N																						
		<input type="checkbox"/> Other	Type: <input type="text" value=""/>																						
Comments, Additional Analyses, or Special Instructions: ONLY PCS RESULTS ABOVE ARE REPORTED. SE-SS-314 ARE REPORTED.																									
Email Results to: REXX@summitenv.com																									
Turnaround Time (TAT) <input type="checkbox"/> 1 Day* <input type="checkbox"/> 2 Days* <input type="checkbox"/> 3 Days* <input type="checkbox"/> 4 Days* <input type="checkbox"/> 5 Days <input checked="" type="checkbox"/> Standard (8-10 business days)																									
Please note: For volatile analyses, a trip blank has been provided in the cooler. If you want the trip blank run and reported please write the trip blank on the COC. Trip Blank analyses will be charged unless other arrangements have been made.																									
*Fee may apply: lab approval required																									
Sampler Name (Print): <input type="text" value="John Cresson"/>																									
Relinquished By Sampler: <input type="text" value="John Cresson"/>																									
Relinquished By: <input type="text" value="John Cresson"/>																									
Relinquished By: <input type="text" value="John Cresson"/>																									
Date: <input type="text" value="2-21"/> Time: <input type="text" value="9:30"/> Received By: <input type="text" value="General Counsel"/>																									
Date: <input type="text" value="2-21"/> Time: <input type="text" value="8:30"/> Received By: <input type="text" value="John Cresson"/>																									
Date: <input type="text" value="2-21"/> Time: <input type="text" value="8:30"/> Received By: <input type="text" value="John Cresson"/>																									



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 74887
CLIENT: Summit
PROJECT: Maine Energy

COOLER NUMBER: 255
NUMBER OF COOLERS: 1

A: PRELIMINARY EXAMINATION:

1. Cooler received by(initials): JB

DATE COOLER RECEIVED/OPENED: 2/22/13

2. Circle one: Hand delivered
(If so, skip 3)

Shipped

Y

N

3. Did cooler come with a shipping slip?

3a. Enter carrier name and airbill number here:

504730

4. Were custody seals on the outside of cooler?

How many & where: _____ Seal Date: _____ Seal Name: _____

Y

N

5. Did the custody seals arrive unbroken and intact upon arrival?

Y

N/A

6. COC#:

7. Were Custody papers filled out properly (ink,signed, legible, project information etc)?

Y

N

8. Were custody papers sealed in a plastic bag?

Y

N

9. Did you sign the COC in the appropriate place?

Y

N

10. Was enough ice used to chill the cooler?

Y N

Temp. of cooler:

40C

B. Log-In: Date samples were logged in:

2/22/13

By: JB

11. Were all bottles sealed in separate plastic bags?

Y

N

12. Did all bottles arrive unbroken and were labels in good condition?

Y

N

13. Were all bottle labels complete(ID,Date,time,etc.)

Y

N

14. Did all bottle labels agree with custody papers?

Y

N

15. Were the correct containers used for the tests indicated:

Y

N

16. Were samples received at the correct pH?

Y

N/A

17. Was sufficient amount of sample sent for the tests indicated?

Y

N

18. Were all samples submitted within holding time?

Y

N

19. Were all containers used within AEL's expiration date?**

Y

N/A

20. Were VOA samples absent of greater than pea-sized bubbles?

Y

N/A

(Note:Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

*If NO, List Sample ID's, Lab #s: _____

When bubbles are present in VOA samples they are labelled from smallest (or no bubbles) to largest. Lab to analyze VOA samples with no bubbles or
smallest bubbles first

20. Laboratory labeling verified by (initials): KEZ

Date: 02/22/13

**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.

Insured Delivery

P.O. Box 1072
Portland, ME 04104

385 Main Street
So. Portland, ME 04106

1-800-698-5035
767-6004.

FAX: 767-7159

E-mail: misraelson@generalcourier.com



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O			
M			

Summit Environmental
640 Main St
Kenison, Me

S H Analytics Lab
I G5 Commerce Way
P O Personville, N.H.
T O

BILLING INFORMATION

QUANTITY	DESCRIPTION	PRICE
1	Jacyn	Time 856 2/22/13

Signature

March 6, 2013

Mr. John Cressey
Summit Environmental
640 Main Street
Lewiston ME 04240

RE: Analytical Results Case Narrative
Analytics # 74946
Maine Energy
Project No: 12-3259.1

Dear Mr. Cressey;

Enclosed please find the analytical results for samples submitted for the above-mentioned project. The attached Cover Page lists the sample IDs, Lab tracking numbers and collection dates for the samples included in this deliverable.

Samples were analyzed for Polychlorinated Biphenyls (PCBs) by EPA Method 8082A.

Unless otherwise noted in the Non-conformance Summary listed below, all of the quality control (QC) criteria including initial calibration, calibration verification, surrogate recovery, holding time and method accuracy/precision for these analyses were within acceptable limits.

This Level II data package has been assembled in the following order:

Case Narrative/Non-Conformance Summary
Sample Log Sheet - Cover Page
PCB Form 1 Data Sheet for Samples
PCB Form 3 MS/MSD (LCS) Recoveries and Blanks
Chromatograms
Chain of Custody (COC) Forms

QC NON CONFORMANCE SUMMARY

Sample Receipt:

No exceptions.

PCBs by EPA Method 8082:

No results were reported below the quantitation limit.

All samples except 74946-1 and 74946-7 required dilution due to concentrations of PCBs that exceeded the calibration range of the instrument or matrix affect.

The MS/MSD analyzed on sample 74946-1-1 did not meet acceptance criteria for PCB 1260 recoveries due to the parent sample having concentrations of PCB 1260 that exceeded the calibration range of the instrument. In addition the MS/MSD pair had three high RPDs. The laboratory control samples (L030413PSOX/LD030413PSOX) were in control for all recoveries but had high RPDs. Results were reported without qualification.

If you have any questions on this data submittal, please do not hesitate to contact me.

Sincerely,
ANALYTICS Environmental Laboratory, LLC



Stephen Knollmeyer
Laboratory Director

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

Report Number: 74946

Revision: Rev. 0

Re: Maine Energy (Project No: 12-3259.1)

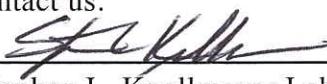
Enclosed are the results of the analyses on your sample(s). Samples were received on 22 February 2013 and analyzed for the tests listed. Samples were received in acceptable condition, with the exceptions noted below or on the chain of custody. These results pertain to samples as received by the laboratory and for the analytical tests requested on the chain of custody. The results reported herein conform to the most current NELAC standards, where applicable, unless otherwise narrated in the body of the report. Please see individual reports for specific methodologies and references.

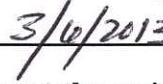
<u>Lab Number</u>	<u>Sample Date</u>	<u>Station Location</u>	<u>Analysis</u>	<u>Comments</u>
74946-1	02/20/13	SE-SB-236 (0-2')	EPA 8082 (PCBs only)	
74946-2	02/20/13	SE-SB-236 (2-4')	EPA 8082 (PCBs only)	
74946-3	02/20/13	SE-SB-236 (4-6')	EPA 8082 (PCBs only)	
74946-4	02/20/13	SE-SB-236 (6-8')	EPA 8082 (PCBs only)	
74946-5	02/20/13	SE-SB-236 (8-10')	EPA 8082 (PCBs only)	
74946-6	02/20/13	SE-SB-236 (10-12')	EPA 8082 (PCBs only)	
74946-7	02/20/13	SE-SB-236 (12-14')	EPA 8082 (PCBs only)	

Sample Receipt Exceptions: None

Analytics Environmental Laboratory is certified by the states of New Hampshire, Maine, Massachusetts, Connecticut, Rhode Island, Virginia, Maryland, and North Carolina, and is accredited by the Department of Defense (DOD) ELAP program. A list of actual certified parameters is available upon request.

If you have any questions on these results, please do not hesitate to contact us.

Authorized signature 
 Stephen L. Knollmeyer Lab. Director

Date 

**This report shall not be reproduced, except in full, without the written
 consent of Analytics Environmental Laboratory, LLC.**

Surrogate Compound Limits

Matrix: Units:	Aqueous % Recovery	Solid % Recovery	Method
Volatile Organic Compounds - Drinking Water			
1,4-Difluorobenzene	70-130		EPA 524.2
Bromofluorobenzene	70-130		
1,2-Dichlorobenzene-d4	70-130		
Volatile Organic Compounds			
1,2-Dichloroethane-d4	70-120	70-120	EPA 624/8260B
Toluene-d8	85-120	85-120	
Bromofluorobenzene	75-120	75-120	
Semi-Volatile Organic Compounds			
2-Fluorophenol	20-110	35-105	EPA 625/8270C
d5-Phenol	15-110	40-100	
d5-nitrobenzene	40-110	35-100	
2-Fluorobiphenyl	50-110	45-105	
2,4,6-Tribromophenol	40-110	40-125	
d14-p-terphenyl	50-130	30-125	
PAH's by SIM			
d5-nitrobenzene	21-110	35-110	EPA 8270C
2-Fluorobiphenyl	36-121	45-105	
d14-p-terphenyl	33-141	30-125	
Pesticides and PCBs			
2,4,5,6-Tetrachloro-m-xylene (TCX)	46-122	40-130	EPA 608/8082
Decachlorobiphenyl (DCB)	40-135	40-130	
Herbicides			
Dichloroacetic acid (DCAA)	30-150	30-150	
Gasoline Range Organics/TPH Gasoline			
Trifluorotoluene TFT (FID)	60-140	60-140	MEDEP 4217/EPA 8015
Bromofluorobenzene (BFB) (FID)	60-140	60-140	
Trifluorotoluene TFT (PID)	60-140	60-140	
Bromofluorobenzene (BFB) (PID)	60-140	60-140	
Diesel Range Organics/TPH Diesel			
m-terphenyl	60-140	60-140	MEDEP 4125/EPA 8015/CT ETPH
Volatile Petroleum Hydrocarbons			
2,5-Dibromotoluene (PID)	70-130	70-130	MADEP VPH May 2004 Rev1.1
2,5-Dibromotoluene (FID)	70-130	70-130	
Extractable Petroleum Hydrocarbons			
1-chloro-octadecane (aliphatic)	40-140	40-140	MADEP EPH May 2004 Rev1.1
o-Terphenyl (aromatic)	40-140	40-140	
2-Fluorobiphenyl (Fractionation)	40-140	40-140	
2-Bromonaphthalene (fractionation)	40-140	40-140	

PCB
DATA SUMMARIES

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (0-2')

Lab Sample ID: 74946-1
Matrix: Solid
Percent Solid: 96
Dilution Factor: 1.0
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit µg/kg	Results µg/kg
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	174

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	92	%
Decachlorobiphenyl	79	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG:

GC Column #1: STX-CLPesticides I

Sample: 74946-1,,A/C

Column ID: 0.25 mm

Data File: M67921.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.0

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	174	166	4.7	

Column to be used to flag RPD values greater than QC limit of 40%

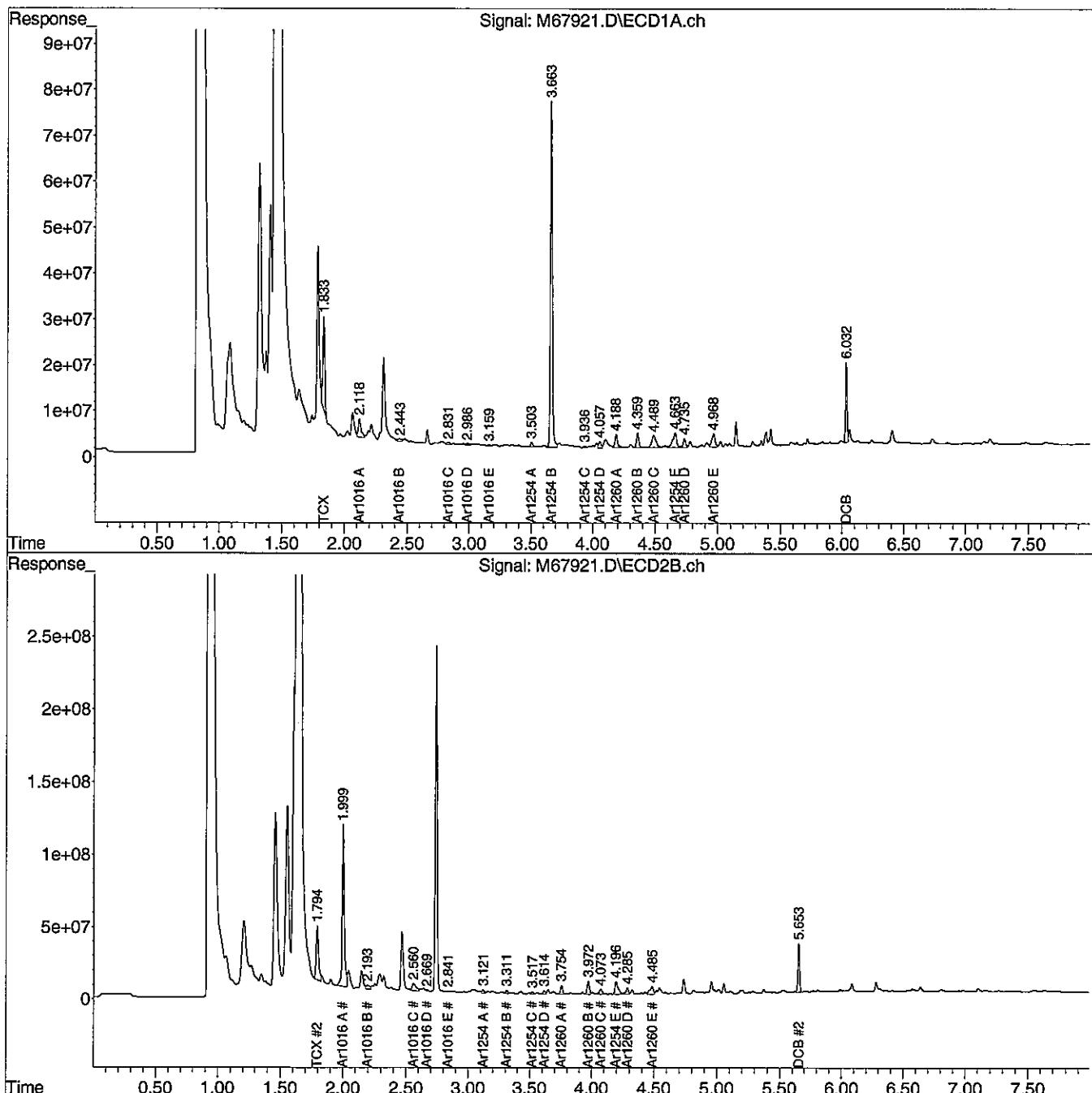
* Values outside QC limits

Comments: _____

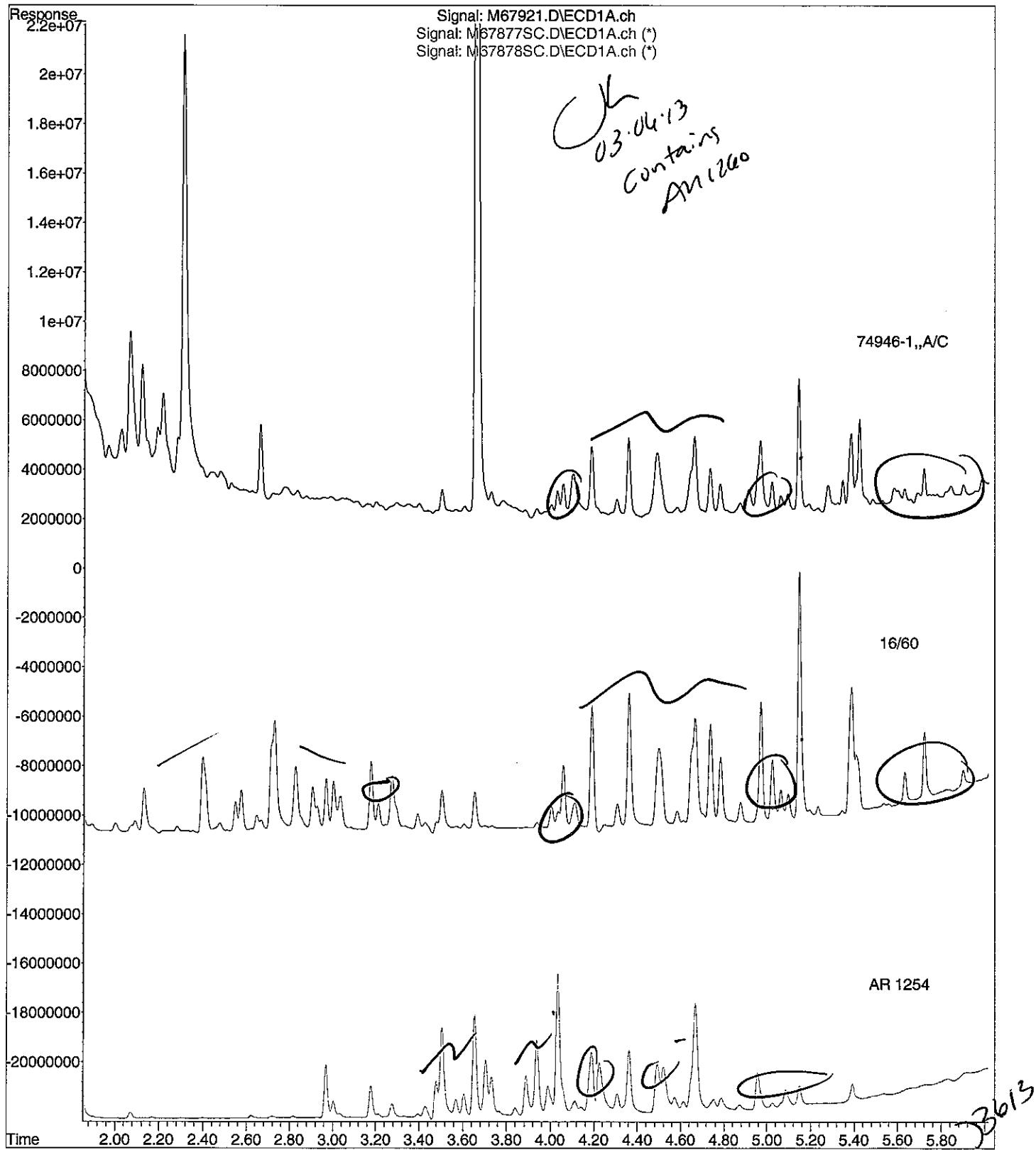
Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67921.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 5:49 pm
 Operator : JK
 Sample : 74946-1, , A/C
 Misc : SOIL
 ALS Vial : 19 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:45:54 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



File : C:\msdchem\1\DATA\030513-M\M67921.D
Operator : JK
Acquired : 5 Mar 2013 5:49 pm using AcqMethod PCB.M
Instrument : Instrument M
Sample Name: 74946-1,,A/C
Misc Info : SOIL
Vial Number: 19



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (2-4')

Lab Sample ID: 74946-2
Matrix: Solid
Percent Solid: 93
Dilution Factor: 2.1
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	69	U
PCB-1221	69	U
PCB-1232	69	U
PCB-1242	69	U
PCB-1248	69	U
PCB-1254	69	U
PCB-1260	69	670

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	112	%
Decachlorobiphenyl	89	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-2,1:2,A/C

Column ID: 0.25 mm

Data File: M67924.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	662	670	1.2	

Column to be used to flag RPD values greater than QC limit of 40%

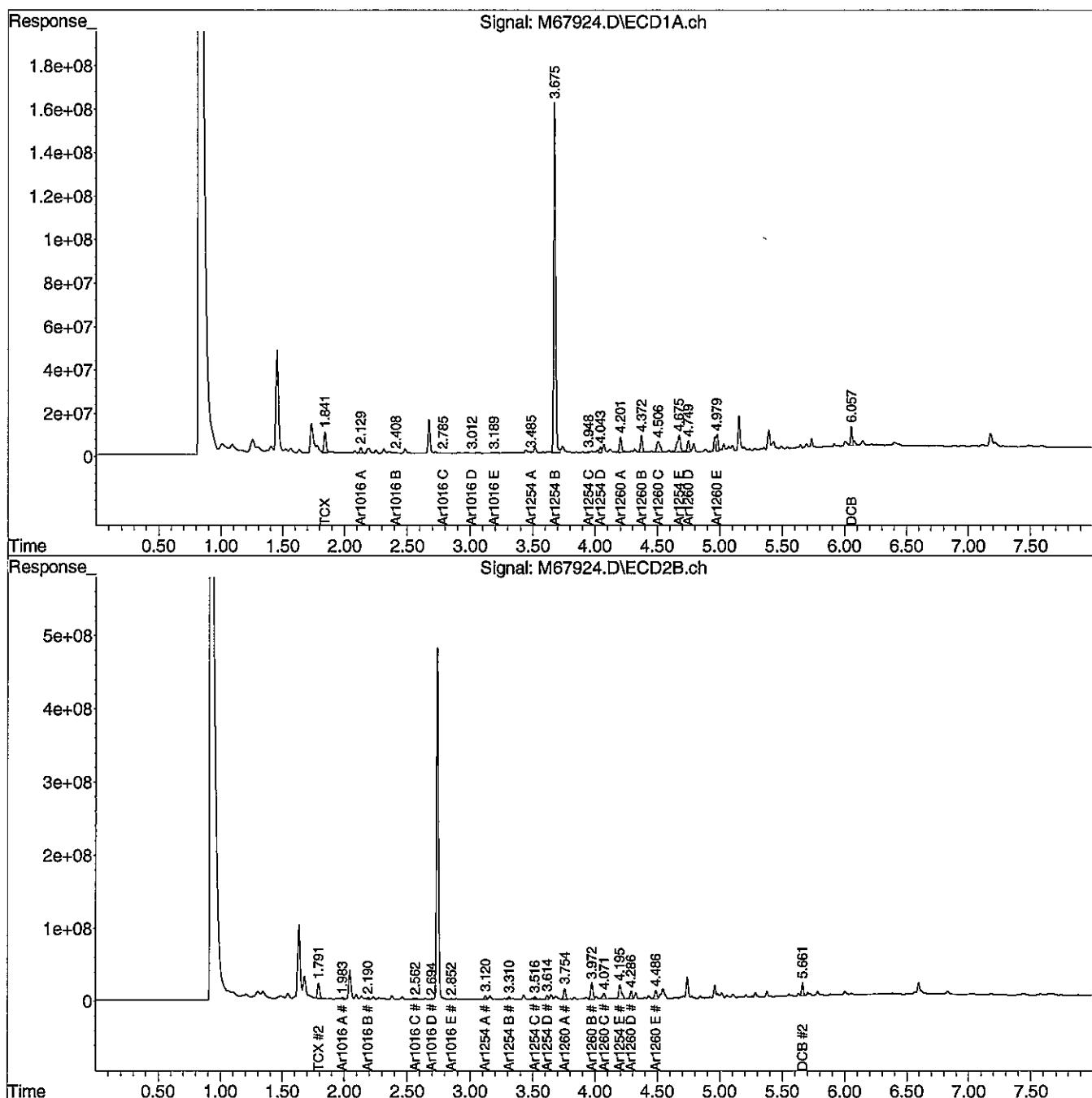
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67924.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 9:15 pm
 Operator : JK
 Sample : 74946-2,1:2,,A/C
 Misc : SOIL
 ALS Vial : 22 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:52:01 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (4-6')

Lab Sample ID: 74946-3
Matrix: Solid
Percent Solid: 89
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	2240

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	124	%	
Decachlorobiphenyl	97	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
--------------	-------------	-----------------------------	---------------------

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-3,1:5,,A/C

Column ID: 0.25 mm

Data File: M67925.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2236	2214	1.0	

Column to be used to flag RPD values greater than QC limit of 40%

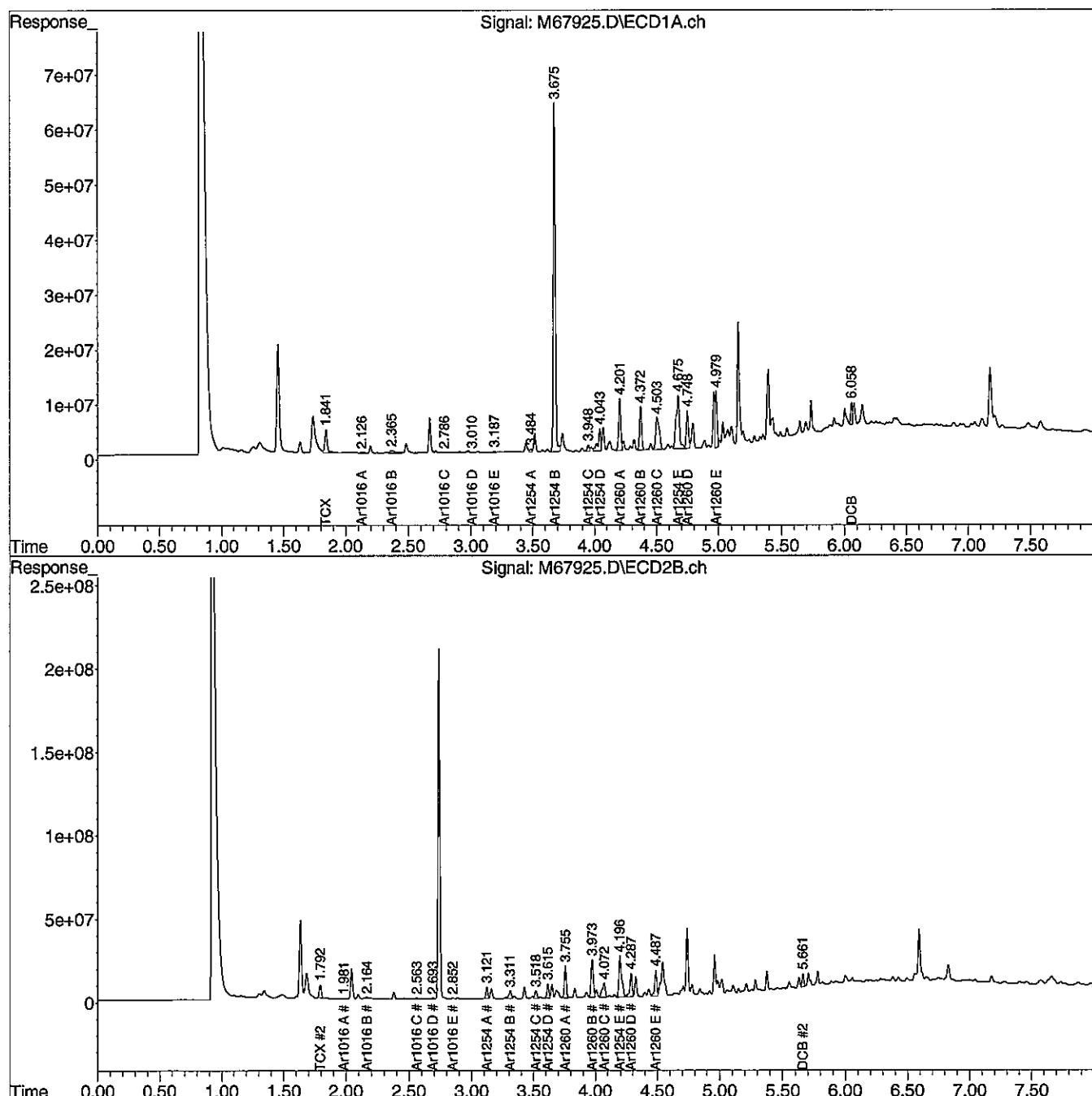
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67925.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 9:44 pm
 Operator : JK
 Sample : 74946-3,1:5,,A/C
 Misc : SOIL
 ALS Vial : 23 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:53:05 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 μ L
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 μ m



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (6-8")

Lab Sample ID: 74946-4
Matrix: Solid
Percent Solid: 91
Dilution Factor: 6
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	198	U
PCB-1221	198	U
PCB-1232	198	U
PCB-1242	198	U
PCB-1248	198	U
PCB-1254	198	U
PCB-1260	198	2120

<u>Surrogate Standard Recovery</u>		
2,4,5,6-Tetrachloro-m-xylene	89	%
Decachlorobiphenyl	77	%

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

PCB
COLUMN RELATIVE PERCENT DIFFERENCE

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-4,1:5,,A/C

Column ID: 0.25 mm

Data File: M67926.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 5.5

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	2121	2062	2.8	

Column to be used to flag RPD values greater than QC limit of 40%

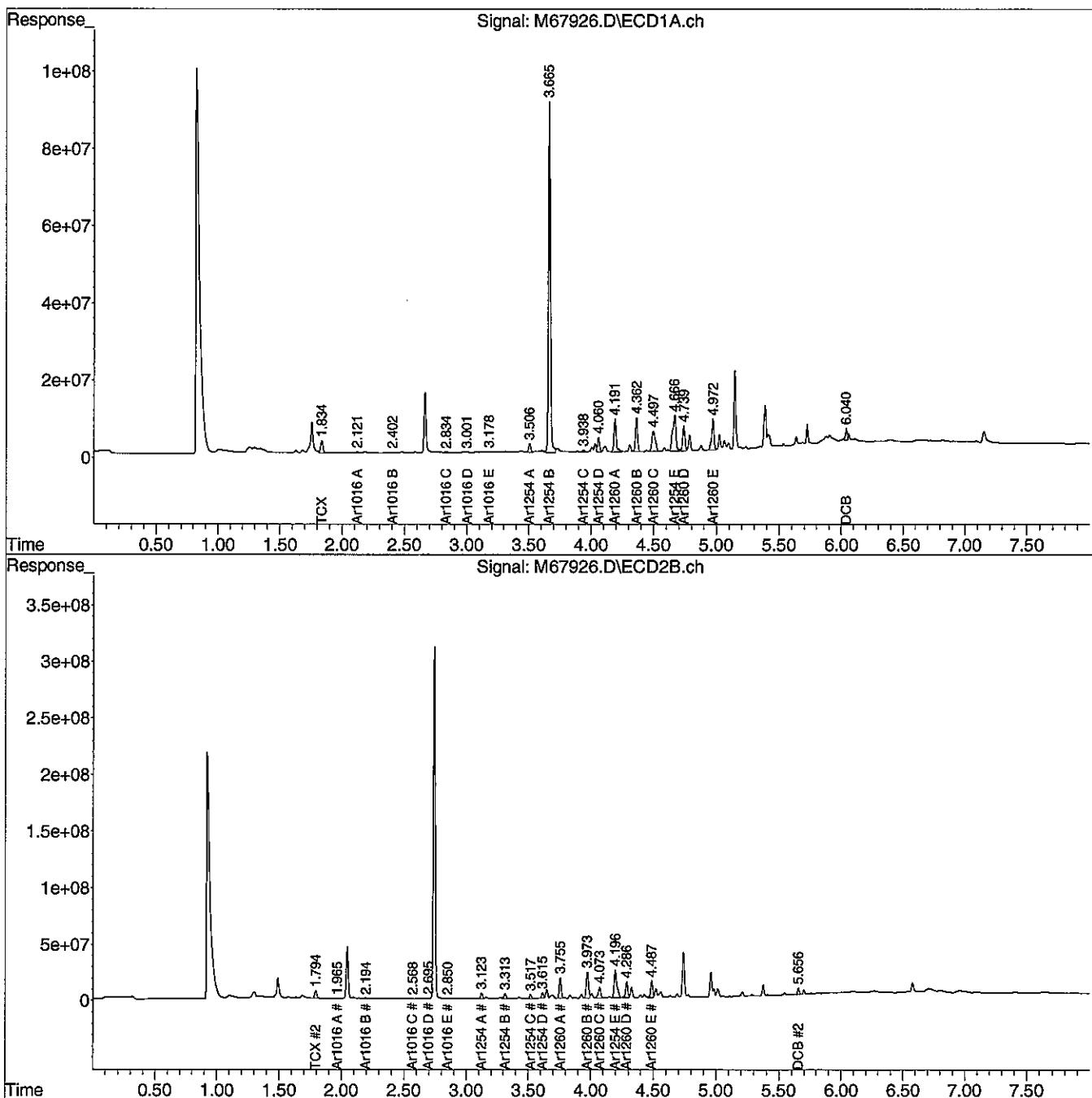
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67926.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 9:54 pm
 Operator : JK
 Sample : 74946-4,1:5,,A/C
 Misc : SOIL
 ALS Vial : 24 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:53:53 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (8-10')

Lab Sample ID: 74946-5
Matrix: Solid
Percent Solid: 90
Dilution Factor: 2.2
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	73	U
PCB-1221	73	U
PCB-1232	73	U
PCB-1242	73	U
PCB-1248	73	U
PCB-1254	73	U
PCB-1260	73	975

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	70	%	
Decachlorobiphenyl	62	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-5,1:2,,A/C

Column ID: 0.25 mm

Data File: M67927.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.2

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	975	948	2.8	

Column to be used to flag RPD values greater than QC limit of 40%

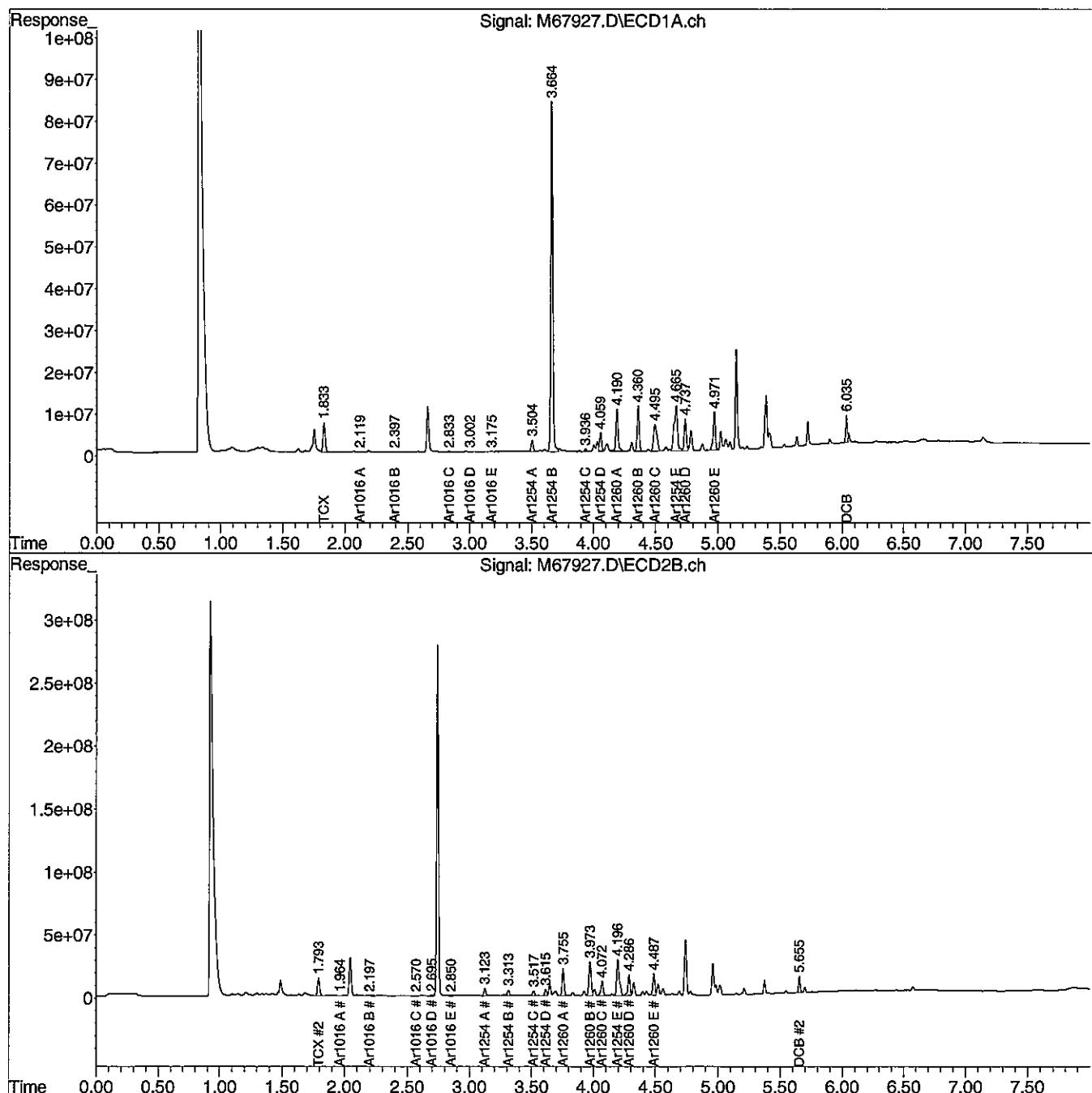
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67927.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 10:04 pm
 Operator : JK
 Sample : 74946-5,1:2,,A/C
 Misc : SOIL
 ALS Vial : 25 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:54:51 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
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 640 Main Street
 Lewiston ME 04240

March 6, 2013
SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (10-12')

Lab Sample ID: 74946-6
Matrix: Solid
Percent Solid: 93
Dilution Factor: 2.1
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	69	U
PCB-1221	69	U
PCB-1232	69	U
PCB-1242	69	U
PCB-1248	69	U
PCB-1254	69	U
PCB-1260	69	735

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	68	%	
Decachlorobiphenyl	64	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-6,1:2,,A/C

Column ID: 0.25 mm

Data File: M67928.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 2.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	735	719	2.2	

Column to be used to flag RPD values greater than QC limit of 40%

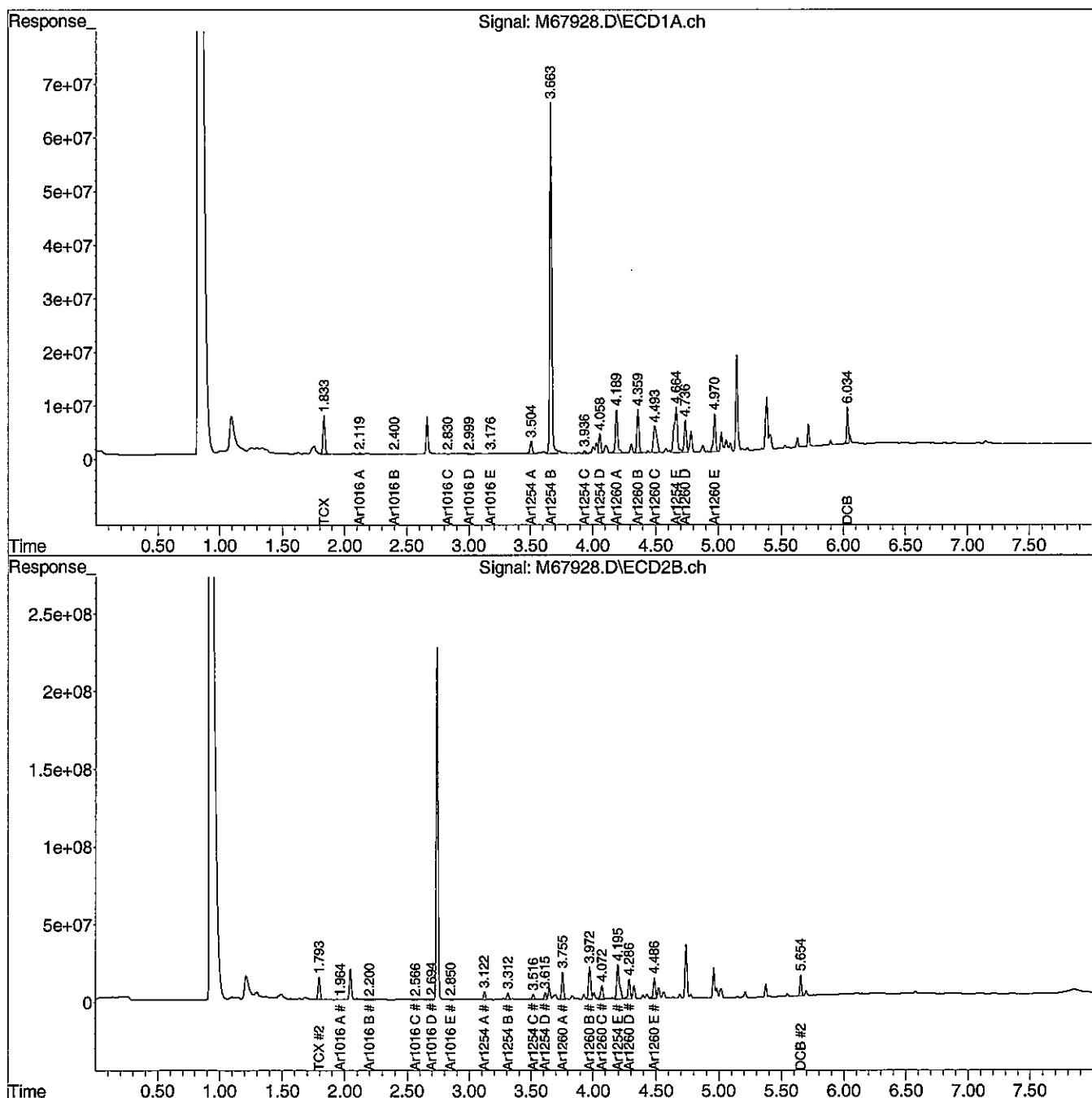
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67928.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 10:14 pm
 Operator : JK
 Sample : 74946-6,1:2,,A/C
 Misc : SOIL
 ALS Vial : 26 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:55:48 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: SE-SB-236 (12-14')

Lab Sample ID: 74946-7
Matrix: Solid
Percent Solid: 88
Dilution Factor: 1.1
Collection Date: 02/20/13
Lab Receipt Date: 02/22/13
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	36	U
PCB-1221	36	U
PCB-1232	36	U
PCB-1242	36	U
PCB-1248	36	U
PCB-1254	36	U
PCB-1260	36	472

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	78	%	
Decachlorobiphenyl	85	%	

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

**PCB
COLUMN RELATIVE PERCENT DIFFERENCE**

Instrument ID: M

SDG: 74946

GC Column #1: STX-CLPesticides I

Sample: 74946-7,,A/C

Column ID: 0.25 mm

Data File: M67929.D

GC Column #2: STX-CLPesticides II

Dilution Factor: 1.1

Column ID: 0.25 mm

COMPOUND	Column #1	Column #2	RPD	#
	SAMPLE RESULT (ug/kg)	SAMPLE RESULT (ug/kg)		
PCB 1260	472	447	5.3	

Column to be used to flag RPD values greater than QC limit of 40%

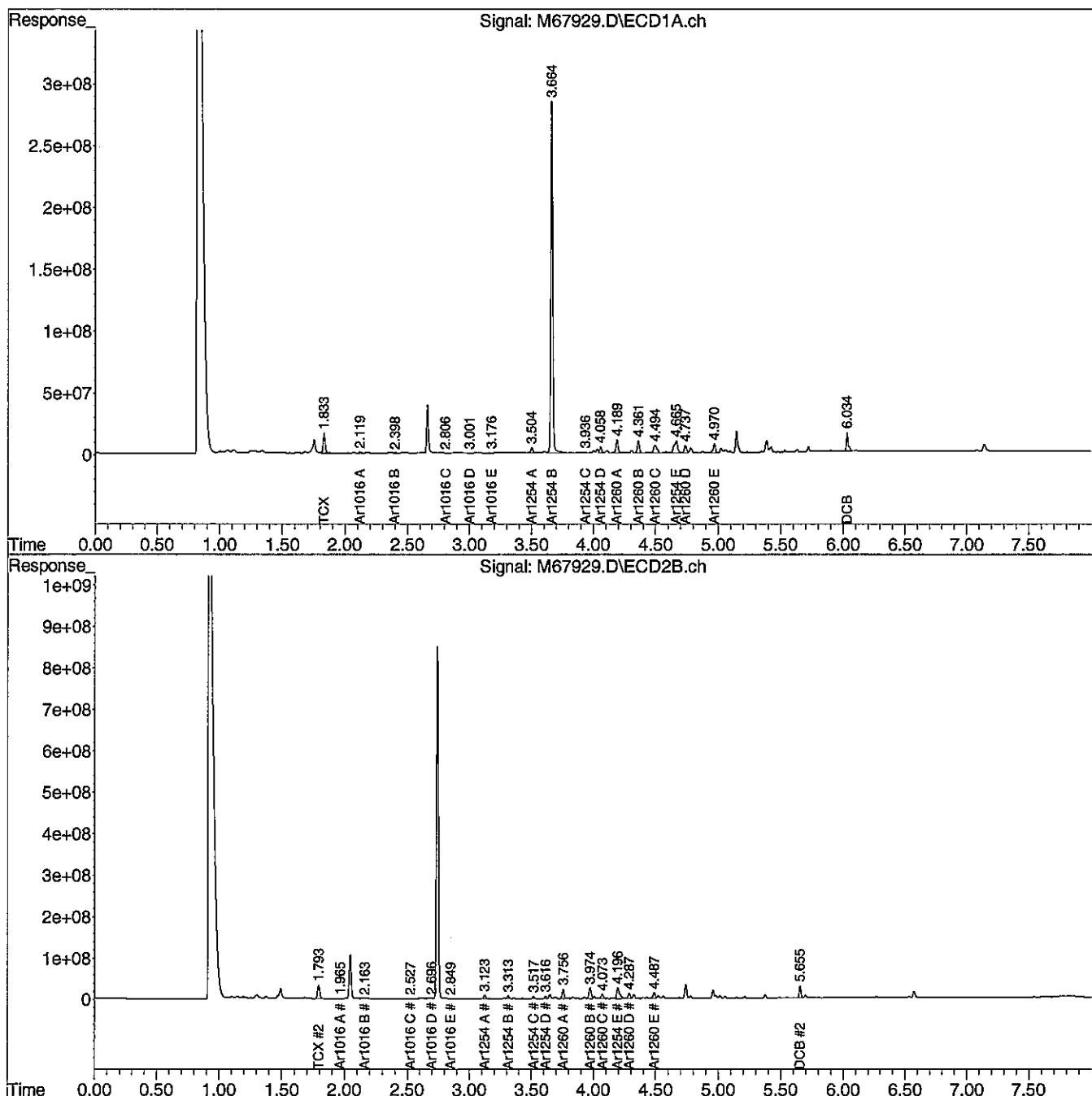
* Values outside QC limits

Comments: _____

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67929.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 10:24 pm
 Operator : JK
 Sample : 74946-7, ,A/C
 Misc : SOIL
 ALS Vial : 27 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:43:25 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:44 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



PCB
QC FORMS

Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
 Project Number: 12-3259.1
 Field Sample ID: Lab QC

Lab Sample ID: B030413PSOX
 Matrix: Soil
 Percent Solid: 100
 Dilution Factor: 1.0
 Collection Date:
 Lab Receipt Date:
 Extraction Date: 03/04/13
 Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g}/\text{kg}$	Results $\mu\text{g}/\text{kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

<u>Surrogate Standard Recovery</u>			
2,4,5,6-Tetrachloro-m-xylene	103	%	
Decachlorobiphenyl	87	%	

U=Undetected	J=Estimated	E=Exceeds Calibration Range	B=Detected in Blank
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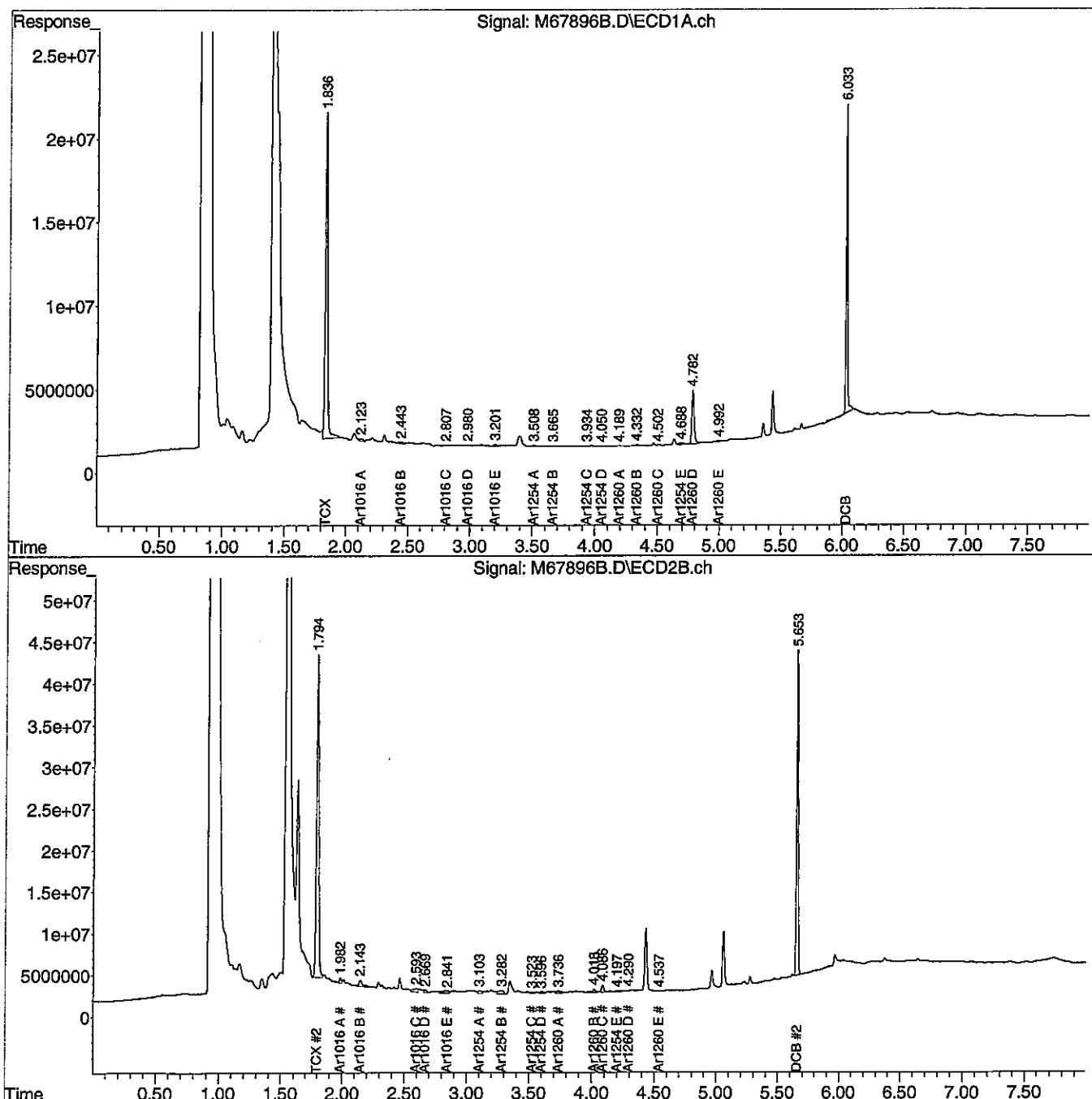
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67896B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 1:37 pm
 Operator : JK
 Sample : B030413PSOX,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 05 14:14:27 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Tue Mar 05 10:44:12 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



Mr. John Cressey
 Summit Environmental Consultants Inc.
 640 Main Street
 Lewiston ME 04240

March 6, 2013

SAMPLE DATA

CLIENT SAMPLE ID

Project Name: Maine Energy
Project Number: 12-3259.1
Field Sample ID: Lab QC

Lab Sample ID: B030413PSOX RR
Matrix: Soil
Percent Solid: 100
Dilution Factor: 1.0
Collection Date:
Lab Receipt Date:
Extraction Date: 03/04/13
Analysis Date: 03/05/13

PCB ANALYTICAL RESULTS

COMPOUND	Quantitation Limit $\mu\text{g/kg}$	Results $\mu\text{g/kg}$
PCB-1016	33	U
PCB-1221	33	U
PCB-1232	33	U
PCB-1242	33	U
PCB-1248	33	U
PCB-1254	33	U
PCB-1260	33	U

Surrogate Standard Recovery

2,4,5,6-Tetrachloro-m-xylene	101	%
Decachlorobiphenyl	86	%

U=Undetected J=Estimated E=Exceeds Calibration Range B=Detected in Blank

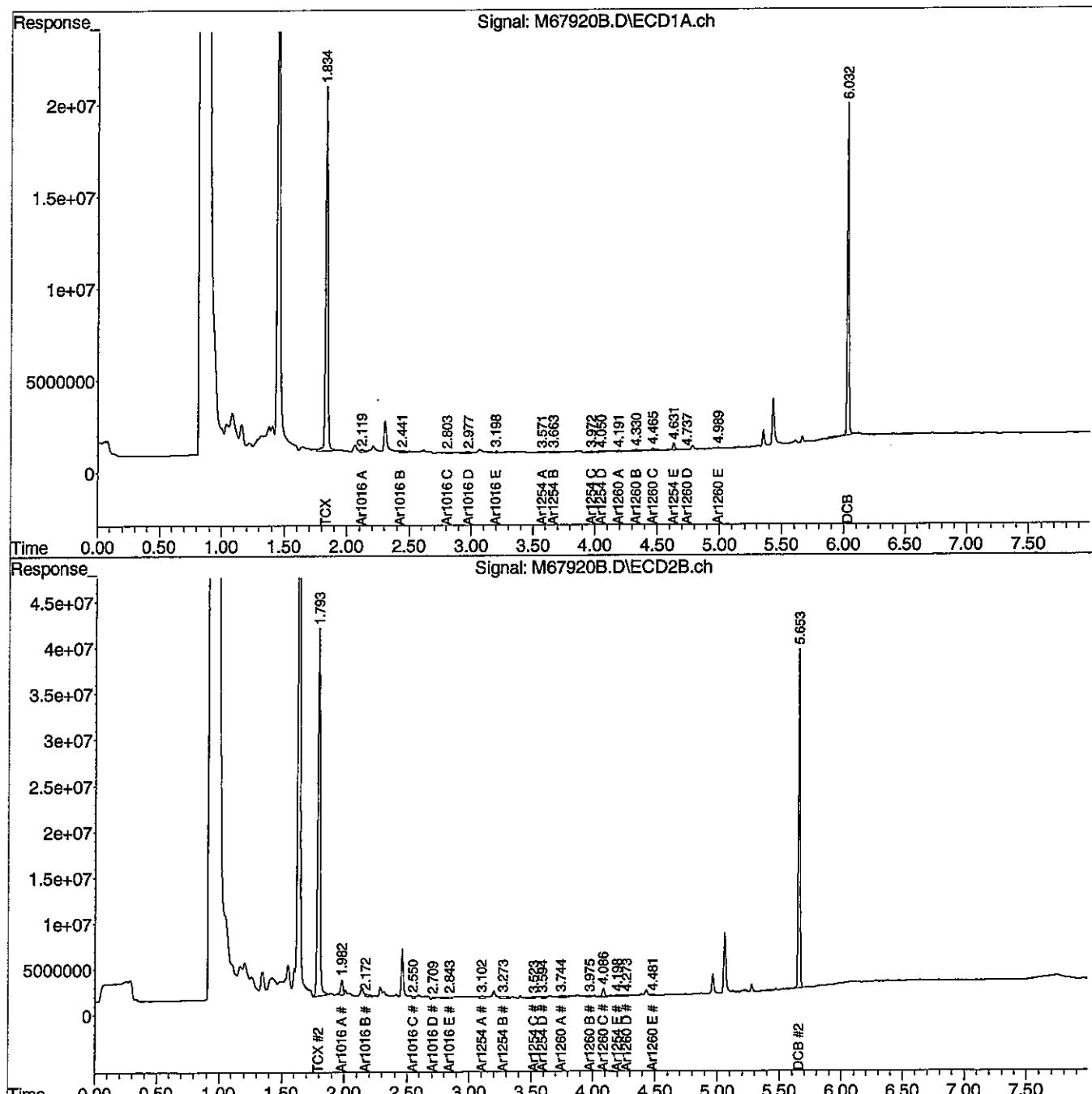
METHODOLOGY: Sample analysis conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 8082A. Sample preparation conducted according to Test Methods for Evaluating Solid Waste, SW-846 Method 3540C. Sample cleanup was conducted according to SW-846 Method 3665A.

COMMENTS: Results are expressed on a dry weight basis.

Data Path : C:\msdchem\1\DATA\030513-M\
 Data File : M67920B.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 5 Mar 2013 5:39 pm
 Operator : JK
 Sample : B030413PSOX,RR2,,A/C
 Misc : SOIL
 ALS Vial : 6 Sample Multiplier: 1

Integration File signal 1: events.e
 Integration File signal 2: events2.e
 Quant Time: Mar 06 10:43:05 2013
 Quant Method : C:\msdchem\1\METHODS\PCB022813.M
 Quant Title : SW-846 METHOD 8082 Aroclor 1016/1260/1254
 QLast Update : Wed Mar 06 10:28:43 2013
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 2 uL
 Signal #1 Phase : STX-CLPPesticides Signal #2 Phase: STX-CLPPesticides
 Signal #1 Info : 30 m x 0.25mm x 0 Signal #2 Info : 30 m x 0.25mm x 0.25 um



PCB SOIL SYSTEM MONITORING COMPOUNDS SUMMARY

Instrument ID: M

GC Column #1: STX-CLPesticides I

SDG: 74946

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits

* Values outside OC limits

D System Monitoring Compound diluted out

**PCB SOIL
SYSTEM MONITORING COMPOUNDS
SUMMARY**

Instrument ID: M
GC Column #1: STX-CLPesticides I
Column ID: 0.25 mm
GC Column #2: STX-CLPesticides II
Column ID: 0.25 mm

SDG: 74946

	Lower Limit	Upper Limit
SMC #1 = TCX	40	130
SMC #2 = DCB	40	130

Column to be used to flag recovery values outside of QC limits
* Values outside QC limits
D System Monitoring Compound diluted out

PCB SOIL
LABORATORY CONTROL SAMPLE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: B030413PSOX,,A/C

Spike: L030413PSOX,,A/C

Spike duplicate: LD030413PSOX,,A/C

COMPOUND	LCS SPIKE	LCSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP		#	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	
PCB 1016	200	200	65	140	30	0	177	89		264	132		39.3	*
PCB 1260	200	200	60	130	30	0	159	79		200	100		23.2	
PCB 1016 #2	200	200	65	140	30	0	155	78		228	114		37.8	*
PCB 1260 #2	200	200	60	130	30	0	163	82		222	111		30.5	*

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

LCS/LCSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

PCB SOIL
MATRIX SPIKE/DUPLICATE
PERCENT RECOVERY

Instrument ID: M

GC Column #1: STX-CLPesticides I

Column ID: 0.25 mm

GC Column #2: STX-CLPesticides II

Column ID: 0.25 mm

SDG:

Non-spiked sample: 74946-1,,A/C

Spike: 74946-1,MS,,A/C

Spike duplicate: 74946-1,MSD,,A/C

COMPOUND	MS SPIKE	MSD SPIKE	LOWER	UPPER	RPD	NON-SPIKE	SPIKE	SPIKE		SPIKE DUP	SPIKE DUP		#	
	ADDED (ug/kg)	ADDED (ug/kg)	LIMIT	LIMIT	LIMIT	RESULT (ug/kg)	RESULT (ug/kg)	% REC	#	RESULT (ug/kg)	% REC	#	RPD	
PCB 1016	203	204	65	140	30	0	204	100		180	88		12.6	
PCB 1260	203	204	60	130	30	174	263	44	*	430	125		48.4	*
PCB 1016 #2	203	204	65	140	30	0	189	93		287	140		41.4	*
PCB 1260 #2	203	204	60	130	30	166	250	42	*	440	134	*	54.9	*

Column to be used to flag recovery and RPD values outside of QC limits

* Values outside QC limits

MS/MSD spike added values have been weight adjusted.

Non-spike result of "0" used in place of "U" to allow calculation of spike recovery.

Comments: _____

CHAIN OF CUSTODIES

Chain Of Custody Form

Chain Of Custody Form



195 Commerce Way, Suite E
Portsmouth, NH 03801
(800) 929-9906

(603) 436-5111
(603) 430-2151 Fax



ANALYTICS SAMPLE RECEIPT CHECKLIST

AEL LAB#: 74946
CLIENT: Summit
PROJECT: Maine Energy

COOLER NUMBER: 255
NUMBER OF COOLERS: 1

A: PRELIMINARY EXAMINATION:

1. Cooler received by (initials): JBDATE COOLER RECEIVED/OPENED: 2/22/132. Circle one: Hand delivered
(If so, skip 3) Shipped

3. Did cooler come with a shipping slip?

 Y N3a. Enter carrier name and airbill number here: 504730

4. Were custody seals on the outside of cooler?

How many & where: _____ Seal Date: _____

 Y

Seal Name: _____

 N

5. Did the custody seals arrive unbroken and intact upon arrival?

 Y N/A

6. COC#:

7. Were Custody papers filled out properly (ink, signed, legible, project information etc)?

 Y N

8. Were custody papers sealed in a plastic bag?

 Y N

9. Did you sign the COC in the appropriate place?

 Y N

10. Was enough ice used to chill the cooler?

 Y NTemp. of cooler: 4°C

B. Log-In: Date samples were logged in:

3/4/13By: JB

11. Were all bottles sealed in separate plastic bags?

 Y N

12. Did all bottles arrive unbroken and were labels in good condition?

 Y N

13. Were all bottle labels complete(ID,Date,time,etc.)

 Y N

14. Did all bottle labels agree with custody papers?

 Y N

15. Were the correct containers used for the tests indicated?

 Y N

16. Were samples received at the correct pH?

 Y N/A

17. Was sufficient amount of sample sent for the tests indicated?

 Y N

18. Were all samples submitted within holding time?

 Y N

19. Were all containers used within AEL's expiration date?**

 Y N/A

20. Were VOA samples absent of greater than pea-sized bubbles?

 Y N/A

(Note: Pea-sized bubbles or smaller are acceptable and are not considered to adversely affect volatiles data.)

*If NO, List Sample ID's, Lab #: _____

When bubbles are present in VOA samples they are labelled from smallest (or no bubbles) to largest. Lab to analyze VOA samples with no bubbles or smallest bubbles first

20. Laboratory labeling verified by (initials): JBDate: 3/4/13

**The expiration date is recommended by Analytics Environmental Laboratory and not the method. Therefore this does not mean that the results are non-compliant.